

# REPORT

## **Source Area 11 2020 Groundwater Report**

**Southeast Rockford Groundwater  
Contamination Superfund Site  
Rockford, Illinois**

968775

Illinois Environmental  
Protection Agency

August 2021

**CDM  
Smith**



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# Section 1

## Introduction

CDM Smith Inc. (CDM Smith) prepared this groundwater monitoring report for the Illinois Environmental Protection Agency (Illinois EPA) to document groundwater quality at Source Area 11 (Area 11) of the Southeast Rockford Groundwater Contamination Superfund Site. The site is located in Rockford, Winnebago County, Illinois, as shown in **Figure 1**.

In addition to documenting the Long-Term Remedial Action (LTRA) for the shallow groundwater component, the data being collected will be used for pre-design evaluations for the soil component remedy that may lead to reexamining the existing remedies.

The first two quarterly rounds of groundwater monitoring in 2020 were conducted in accordance with the Quality Assurance Project Plan (QAPP) Addendum for Area 11 Long Term Remedial Action Monitoring and the Sampling and Analysis Plan (SAP), dated March 19, 2014 and prepared by CDM Smith. The QAPP and SAP were updated per letters from CDM Smith to the U.S. Environmental Protection Agency (USEPA) that were dated May 15, 2019 and November 4, 2019. The third and fourth quarterly sampling rounds in 2020 were conducted in accordance with the Final QAPP for Source Area 11 Long Term Remedial Action Groundwater Monitoring and the SAP, dated August 2020 prepared by CDM Smith.

The current groundwater monitoring network includes 19 monitoring wells, 9 of which are sampled during groundwater sampling events with the remaining 10 used only for water level measurements. The monitoring well network is shown in more detail in **Figure 2**

The report focuses on the methods and procedures used during the 2020 quarterly monitoring events, presents the data for the groundwater elevation measurements and quarterly analytical results, and summarizes information from monitoring events conducted from 2011 to 2020. This report discusses the current contaminant migration scenario and updated site conceptual model.

### 1.1 Area 11 Background Information

Area 11 is located on the northeast corner of Harrison Avenue and Eleventh Street in Rockford, Illinois. The site consists of mixed greenspace and industrial property. The Accurate Metals – Illinois (AMI) facility (formerly Rohr Manufacturing and Rockwell Graphics Systems) is located to the north, and greenspace owned by the City of Rockford (formerly Villa Di Roma restaurant and Rockford Varnish) is located along the southern end of the property. The greenspace portion was purchased by the City of Rockford for use as a storage and laydown area during reconstruction of Harrison Avenue that occurred between 2016 and early 2018.

Potential contaminant sources include eight aboveground storage tanks (AST) that were located east of the former Rockford Varnish facility, leaking tanks, above ground pipes, and a bunker reportedly used by Rockford Varnish located near a former railroad spur. The ASTs were removed sometime between July 2003 and April 2005 based on Google Earth Pro historical imagery. The specific chemicals stored in individual tanks are not known. **Figure 3-23** from the “Final Remedial Investigation Report for the Southeast Rockford Source Control Operable Unit” dated July 25, 2000, shows the historical site features. Following the City’s purchase of the property in 2014, all of the asphalt and above ground structures (and some

shallow subsurface structures), were removed prior to use during the Harrison Avenue construction that started in 2016. Some concrete foundation walls in the vicinity of the former buildings and ASTs are known to remain, and it is generally believed that most other subsurface structures remain.

The geologic stratigraphy at Area 11 is fine- to medium-grained sand down to about 30 feet below ground surface (bgs), followed by medium- to coarse-grained sand with gravel down to about 75 feet bgs. Below this is a silt and clay layer believed to be around 10 to 15 feet thick. The depth to groundwater is approximately 30 feet bgs and varies seasonally.

Groundwater in the unconsolidated material at Area 11 enters the eastern edge of the site flowing in a northwesterly direction before eventually turning west, and then west-southwest as it exits the site's western boundary. Further downgradient, flow directly to the southwest has also been observed.

This southwest flow has also been documented at Area 9/10. This gradual shift in groundwater flow from the northwest to the southwest in the vicinity of Area 11 is responsible for the "banana" shape of the historic groundwater contaminant plume. This plume was documented by CDM Smith during the remedial investigation phases and is critical to understanding contaminant migration patterns in groundwater at Area 11.

In accordance with the Source Control Operable Unit Record of Decision (ROD) dated June 11, 2002, the Area 11 remedy selected for contaminated soil is soil vapor extraction (SVE). The remedy for "leachate" (i.e., shallow, contaminated groundwater) is no action, with groundwater monitoring and institutional controls. The ROD did not propose a remedial alternative for the treatment of leachate on-site because modeling indicated that groundwater would meet standards by the time it exited the source area. However, the ROD indicates that an air sparging component can be added to the remediation system if an improvement in groundwater quality is not observed.

Contaminants of concern (COC) listed in the ROD include benzene, ethylbenzene, methylene chloride, toluene, trichloroethene, and xylenes. However, based on the magnitude of the remediation goal (RG) exceedances in groundwater samples collected since 2008, ethylbenzene, toluene, and xylenes (ETX) are generally considered to be the primary COCs.

Three rounds of pre-design investigation activities were conducted between 2007 and 2018. The first two rounds of investigation in 2009 and 2013 occurred prior to the Harrison Avenue construction conducted by the City of Rockford in 2018. One common objective was to identify and characterize the source material locations in the vadose zone (i.e., where waste material was deposited) that are the targets of the SVE soil component remedy. However, the precise locations of the vadose contamination have not been located for various reasons including site access issues, buried debris, and the assumed small footprints of the source material. The number of individual sources present at Area 11 is not known but it is believed that at least several exist based on groundwater results. The Phase II Pre-Design Technical Memorandum dated September 10, 2013 contains a comprehensive discussion of the nature and extent of groundwater contamination and the possible source locations at Area 11.

Pre-design objectives that were successfully achieved included defining the extent of groundwater contamination at, and downgradient of Area 11 for the leachate component remedy. As part of the pre-design activities, quarterly groundwater sampling was conducted (with several interruptions from 2011 through 2013), before changing to semiannual

sampling in August 2014 for the start of the leachate component long term remedial action (LTRA). These events have resulted in the soil component of the selected remedy remaining in the remedial design (RD) phase while the leachate component has progressed into LTRA.

From 2015 through 2018, semiannual monitoring was not possible due to various factors. For example, in 2015, only one round of groundwater sampling was performed due to contract issues and in 2016, 2017 and 2018, sampling activities were impacted by the Harrison Avenue construction, allowing only one round of groundwater sampling to be completed. Semiannual sampling resumed in 2019 following completion of the Harrison construction.

The third round of pre-design field activities was conducted in October 2018. This phase of work was conducted after the area adjacent to Harrison Avenue had been cleared of buildings, structures, pavement, and road construction debris. The purpose of the activities was to locate and characterize contaminant source material in the Area 11 vadose zone after the removal of obstructions that impacted previous investigations. This phase of work was narrowly focused on areas upgradient of highly contaminated groundwater, and downgradient of suspected point sources of contamination that had become accessible due to the completion of the construction activities. The planned activities included two trenches and one shallow test pit to be excavated followed by direct push soil and groundwater sampling. Details of the activities are documented in the Technical Memorandum dated December 31, 2018. This third phase of pre-design work was not successful in identifying the location of contaminant source material at Area 11, however, the removal of much of the impervious surfaces in that area does appear to have influenced the concentrations of contaminants in groundwater, as a decrease in concentrations has been observed since the completion of the construction. During the field activities, monitoring well MW-007 was installed at a location immediately downgradient of Area 11.

Based on the groundwater data collected from 2017 through 2019, it was determined that while contaminant concentrations have decreased within Area 11, in the areas where barriers to infiltration have been removed, contaminants (primarily ETX) are now migrating downgradient at levels above the remediation goals. This contamination continues to attenuate rapidly with distance; however, this is a significant difference from what is shown in the basic conceptual site model (CSM). It was recommended that quarterly groundwater monitoring be conducted for a minimum of two years to more fully evaluate the revised CSM for Area 11. In addition to continuing to monitor for volatile organic compounds and 1,4-dioxane, the number of water level measurement locations was increased to enhance the understanding of the hydrogeology. These data provide input to refine the remedial decisions for leachate and soils within Area 11.

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## Section 2

# Field and Analytical Activities

The current groundwater monitoring network includes 19 monitoring wells, as shown in **Figure 2**. Depth to groundwater measurements and groundwater samples were collected from nine monitoring wells and depth to groundwater measurements only were collected from 10 monitoring wells. **Table 1** provides a summary of the groundwater monitoring sampling dates and wells sampled for the 2020 events. The first two quarterly events were performed under the previous QAPP and included 12 monitoring well locations for depth to groundwater measurements. Eight of those well locations were also used for groundwater sample collection. **Table 2** provides the updated monitoring well details.

## 2.1 Groundwater Elevations

Depth to groundwater measurements were collected manually at each well prior to purging and sample collection, except for the September 2020 event, where applicable. Prior to the September 2020 event, three water level only monitoring wells were inaccessible and this round of water measurements was subsequently collected in October 2020.

An electronic water level indicator was used and decontaminated before and after each use. Potentiometric surface maps were prepared from the groundwater elevation data collected during the October and November quarterly water level events in 2020, using data from the 18 monitoring wells screened in the upper portion of the unconsolidated materials (**Figures 3 and 4**). With the larger set of monitoring points during October and November, these events present a more accurate picture of groundwater flow patterns. The groundwater elevation data used to compile these maps is provided in **Table 3**.

## 2.2 Sampling Methods

The Area 11 monitoring wells were each purged using a submersible pump and pump controller capable of operating at low flow purging rates. All wells were purged and sampled in general accordance with the applicable SAP.

For all wells sampled, field measurements of pH, temperature, specific conductance, dissolved oxygen, turbidity, and oxidation-reduction potential were monitored with a flow-through multiparameter probe to identify the point stabilization was observed during purging. Parameter readings were recorded at 5-minute intervals and purging continued until the field parameters were observed to be within stable range for three consecutive readings. The stabilization requirements are provided as follows:

- pH:  $\pm 0.25$  standard units
- Dissolved oxygen:  $\pm 10$  percent
- Specific conductance:  $\pm 50$  millisiemens per centimeter (mS/cm) centimeter
- Turbidity: less than 5 nephelometric turbidity units (NTU) or  $\pm 10$  percent
- Temperature:  $\pm 0.5$  C°

- Oxidation-reduction potential:  $\pm 10$  millivolts (mV)

Final readings taken prior to sampling are provided in **Table 4**, and original data sheets listing all readings recorded during purging are provided in **Appendix A**.

Quality control samples specified in the QAPP for each of the groundwater sampling events included one field duplicate per 10 (or fewer) investigative samples, one field blank per 10 (or fewer) investigative samples collected using non-dedicated equipment, one trip blank for each cooler shipped containing aqueous samples for VOC analysis and 1,4-dioxane analysis by Region 5 Analytical Service Branch (ASB) laboratory, and one matrix spike/matrix spike duplicate (MS/MSD) per 20 (or fewer) samples.

The field duplicate frequency was met for all parameter groups for all four quarterly events. The field blank collection frequency was met for 1,4-dioxane and VOCs for all four quarterly events, however field blank collection frequency was not met for the attenuation parameters for the first two quarterly events. Even though the frequency criteria were not met for anions, methane and alkalinity, data quality objectives are not compromised as these analytes are not constituents of concern and the field blanks that were collected for them did not have high detections of these analytes. Field blank contamination actions are discussed in the individual validation reports. A trip blank was sent with each cooler containing samples for VOC or 1,4-dioxane analysis.

Field instruments were calibrated daily to the appropriate standards, in accordance with the SAP. The field samples collected for dissolved ferrous iron were run through a filter attached to the sample tubing and analyzed in the field with a field test kit. New or dedicated sample tubing was used for each discrete sampling location. The groundwater samples selected for laboratory analysis were collected directly from the pump discharge tubing into pre-preserved sample containers. The sample containers were provided by a commercial laboratory.

## 2.3 Analytical Methods and Laboratories

Groundwater samples for 2020 were analyzed for Target Compound List (TCL) VOCs by U.S. EPA Contract Laboratory Program (CLP) laboratories under Statement of Work (SOW) SOM02.4 or VOCs by Region 5 ASB laboratory using Standard Operating Procedure (SOP) MS023. Groundwater sample analysis of 1,4-dioxane was performed by the ASB laboratory in accordance with ASB SOP MS035 for low-level 1,4-dioxane, or by CLP. The U.S. EPA CLP used Chemtech Consulting Group and Pace Analytical Services laboratories for organic sample analyses. Tech Law Inc. (Tech Law) Environmental Services Assistance Team (ESAT), provided services to Region 5 ASB and STAT Analysis Corporation (STAT), Chicago, Illinois and Eurofins TestAmerica, Savannah, Georgia provided anions, alkalinity, and methane analyses. Field analysis of dissolved ferrous iron was performed in accordance with HACH Method 8146.

## 2.4 Data Evaluation and Usability

A data evaluation/validation review was conducted on the analytical data for the four 2020 quarterly groundwater monitoring events. Quality assurance objectives for measurement data are expressed in terms of precision, accuracy, representativeness, comparability, completeness, and sensitivity (PARCCS). The PARCCS parameters characterize the quality of the data and are called data quality indicators (DQIs). The DQIs provide a mechanism for

ongoing quality control (QC) and evaluating and measuring data quality throughout the project. The measurement performance criteria are outlined in the 2014 QAPP/SAP, modified per letters from CDM Smith to U.S. EPA dated May 15 and November 4, 2019, and the August 2020 QAPP/SAP update.

Reviewing the collected data is necessary to determine if data measurement objectives established in the QAPP were met. In general, the following data measurement objectives were considered:

- Achievement of analytical method and reporting limit requirements
- Adherence to and achievement of appropriate laboratory analytical and field QC requirements
- Achievement of required measurement performance criteria for DQIs (the PARCCS parameters)
- Adherence to sampling and sample handling procedures
- Adherence to the sampling design and deviations documented on field change notifications, if required

Data verification, data validation, and data assessment were used to verify adherence to the QAPP procedures and requirements and achievement of the measurement performance criteria of the PARCCS parameters. These assessments were used to reconcile the planned objectives detailed in the QAPP against the investigation results. The outputs serve to verify that the collected data are of sufficient quality to support their intended use.

There were 25 sample delivery groups from the CLP laboratories, Tech Law, STAT and Eurofins. Validation was performed following the Stage 2B validation requirements, EPA's current National Functional Guidelines, current CLP SOWs, and the Region V Organic CLP validation SOP 83074-8-33-601-SO-1143.R1. In accordance with the QAPP, the Tech Law, STAT and Eurofins data were validated by CDM Smith at a Stage 2B Validation/Verification level. The CLP data was validated by the USEPA. CDM Smith reviewed the CLP validation reports and verified the sample results and qualifiers.

The detailed data evaluation/validation discussion is provided as a preface to the laboratory data reports in **Appendix B**. Some analytes were qualified as estimated (J), estimated biased high (J+) or biased low (J-) and/or non-detect (U) or estimated non-detect (UJ), based on validation criteria. Specific details on qualifications are provided in the individual data validation reports.

All field duplicate relative percent difference (RPD) results were within appropriate criteria except for field duplicate pair A11-MW007-201201/A11-MW007-201201-D. Sample results for isopropylbenzene, n-propylbenzene, sec-butylbenzene, 1,3,5-trimethylbenzene, benzene, naphthalene, and n-butylbenzene were qualified as estimated (J/UJ) based on RPD criteria or absolute difference criteria. For this field duplicate pair, the RPD criteria was met for the anions, alkalinity, and methane results.

A review was conducted on the VOC analyses as the sample concentrations vary enough to be suspect based on past sampling results. The COCs were evaluated for possible sample label issues and a comparison was done between samples that were collected on the same day. The

COCs indicated no mislabeled samples and the sample comparison showed no other sample that had similar results comparable to A11-MW007 or A11-MW007-D. A review of the raw laboratory data and chromatograms showed no obvious system errors. The VOC sample results that did not meet RPD criteria are estimated following data validation guidance and there is the potential a sample mix up may have occurred in the field or the laboratory. The results should be used with caution and future sampling events at these locations will be conducted to evaluate the variable sample results.

In summary, all the validated and reviewed data are suitable for their intended use for site characterization. No data were rejected for the 2020 sampling events. Sample results that were qualified as estimated are usable for project decisions. Results that have been rejected from previous sampling years are not usable for project decisions. The laboratory and validation qualifiers are provided in the data tables referenced in **Section 3**.



## Section 3

### Results

This section presents the results of the four quarterly groundwater sampling events in 2020. The Area 11 monitoring wells include two wells upgradient (MW-001 and MW-130A), four wells within Area 11 (MW-002, MW-003, MW-004A, and MW-004B) and three wells downgradient of Area 11 (MW-005, MW-006, and MW-007). Downgradient well MW-007 was installed in 2018. Upgradient well MW-130A was added to the well network for the September and December sampling events. Two of the wells within Area 11 (MW-004A and MW-004B) are adjacent, with MW-004A screened in the shallow portion of the aquifer, and MW-004B screened 45 feet lower in order to monitor the vertical extent of the groundwater contamination within the source area. Because the primary COCs are less dense than water, MW-004B also provides an indication of diffused sitewide contamination.

The monitoring well sample concentrations for VOCs were compared to the remediation goals that were established in the ROD. The sample concentrations for 1,4-dioxane were compared to the Illinois EPA groundwater quality Class I potable groundwater standard provided in Illinois Administrative Code (IAC) 35, 620.410.

#### 3.1 Hydraulic Results

Groundwater elevation measurements were collected prior to the start of each quarterly sampling event. The dates of data collection and the water elevations measured for the 2020 groundwater monitoring events are presented in **Table 3**. Potentiometric surface maps are presented for the last two quarterly events in **Figures 3 and 4**. These events were selected because they present a more complete picture of groundwater flow patterns. Groundwater in the unconsolidated material enters the eastern edge of Area 11, flowing in a northwesterly direction before turning west-southwest as it exits Area 11 along the western boundary. Due to this gradual shift in groundwater flow direction, two gradients that run perpendicular to groundwater flow are estimated for Area 11. From the eastern boundary to the shift in direction near Harrison Avenue, groundwater gradients are estimated using elevation data from MW-32 (as the upgradient location), and MW-004A (as the downgradient location). After the flow direction shifts, the gradients are estimated using elevation data from MW-0007 (as the upgradient location), and MW-126A (as the downgradient location).

The third 2020 quarterly groundwater elevations were measured on October 15, 2020. The groundwater flow direction was measured predominantly to the west, as shown in **Figure 3**, with the bend in flow direction occurring near Harrison Avenue. This quarterly event includes water level measurements from 18 monitoring wells, and the groundwater gradient from MW-32 to MW-004A was approximately 0.003971 feet/foot. The gradient from MW-007 to MW-126A was approximately 0.001411 feet/foot, which is consistent with the historical average groundwater gradients.

The fourth 2020 semiannual groundwater elevations were measured on November 30, 2020. The groundwater flow direction was measured predominantly to the west, as shown in **Figure 4**, with the bend in flow direction occurring near Harrison Avenue. This quarterly event includes water level measurements from 18 monitoring wells. The groundwater gradient from MW-32 to MW-004A was approximately 0.003911 feet/foot, which is consistent

with the October 2020 gradient. The gradient from MW-007 to MW-126A was approximately 0.001360 feet/feet, which is consistent with the historical average groundwater gradients.

## 3.2 Laboratory Analytical Results for VOCs

The laboratory analytical results for the monitoring wells were compared to the remediation goals (RG) from the OU3 ROD and to the “Groundwater Quality Standards for Class I: Potable Resource Groundwater, Title 35, Illinois Administrative Code (35 IAC), Section 620.410”.

**Table 5** summarizes VOCs detected during the 2020 quarterly groundwater monitoring events. **Table 6** summarizes the VOCs that have been detected in at least one sample collected during the baseline, or subsequent, monitoring events for each well. Detected compounds are shown in bold type, and compounds exceeding their RG are shaded. For sample locations where an investigative and field duplicate sample were analyzed, if one of the two (or both) exceeds an RG for a parameter, the location is described as exceeding RGs in the results discussion. Complete analytical results and data validation reports are provided in **Appendix B**.

The groundwater monitoring investigative samples and associated QC samples were analyzed and the data validated as described in **Sections 2.3** and **2.4**. Due to the differences in analytical methods for VOCs used by the different laboratories (CLP and ASB), the VOC parameter lists analyzed by each of the different laboratories are slightly different. In **Tables 5 and 6**, any parameters not analyzed for a particular sampling event because of laboratory assignment are designated with “NA” for not analyzed. MW-130A was added to the monitoring well network in August 2020.

A recurring issue is that high concentrations of ETX compounds (typically toluene) in samples collected from MW-002 and MW-004A routinely require dilution which results in elevated detection limits for all other compounds that are greater than their respective RG. For the 2020 sampling year, this primarily occurred for the compounds with the lower RGs. Because of this “masking” effect, it is not possible to determine if other compounds are present above their RGs in the samples.

### 3.2.1 First Quarter (March 2020) Volatile Organic Compounds Exceeding Remediation Goals

During the March 2020 groundwater sampling event, eight Area 11 wells were sampled. MW-002 continues to have the most compounds exceeding RGs at Area 11. Ethyl benzene was detected at 10 times the RG of 700 µg/L, toluene was detected at 79 times the RG of 1,000 µg/L, and xylenes were detected at two- and one-half times the RG of 10,000 µg/L. Well M-3 contained ethyl benzene at two times the RG, and xylenes slightly over the RG. Well MW-004A contained toluene at 45 times the RG. Ethyl benzene was detected in both the sample and field duplicate at MW-007 slightly over the RG. The compound 1,4-dioxane was detected at slightly over the RG of 7.7 µg/L in MW-003 and MW-004B (**Table 5**).

### 3.2.2 Second Quarter (June 2020) Volatile Organic Compounds Exceeding Remediation Goals

Eight Area 11 monitoring wells were sampled during the June 2020 sampling event. The most contaminated well, MW-002, had ethyl benzene at 9 times the RG, toluene at 68 times the RG and xylenes at 2.5 times the RG. Additionally, vinyl chloride was detected at two times the RG in MW-002. The sample from MW-004A contained toluene at 52 times the RG and tetrachloroethene slightly above the RG. The sample and field duplicate from MW-007

contained ethyl benzene slightly over the RG and MW-005 contained bromodichloromethane at two times the RG of 0.2 µg/L. The compound 1,4-dioxane was detected at slightly over the RG in MW-003, MW-004B and MW-005 and at two times the RG in MW-001 (**Table 5**).

### 3.2.3 Third Quarter (September 2020) Volatile Organic Compounds Exceeding Remediation Goals

For the third quarter monitoring event, nine wells were sampled because MW-130A was added to the monitoring well network in August 2020. In well MW-002, ethyl benzene was detected at 12 times the RG, toluene was detected at 39 times the RG, and xylenes were detected at 3 times the RG. Well MW-004A contained toluene at 43 times the RG. The sample and field duplicate from MW-007 contained ethyl benzene that was three and a half times the RG. The compound 1,4-dioxane was detected at slightly over the RG in MW-004B, MW-005 and MW-006 (**Table 5**).

### 3.2.4 Fourth Quarter (December 2020) Volatile Organic Compounds Exceeding Remediation Goals

All nine Area 11 monitoring wells were sampled during the December 2020 quarterly sampling event. During all four quarters of the 2020 sampling, MW-002 continued to have the most compounds exceeding RGs at Area 11. For the fourth quarter, ethyl benzene was detected at 15 times the RG, toluene was detected at 33 times the RG, and xylenes were detected at four times the RG. MW-004A contained toluene at 34 times the RG. The MW-007 sample and field duplicate contained ethyl benzene at four times the RG, and benzene was detected in the field duplicate only at eight times the RG of 5 µg/L. The compound 1,4-dioxane was not detected above the RG in any well for this sampling round. (**Table 5**).

### 3.2.5 Comprehensive Compounds Exceeding Remediation Goals

The concentrations of different contaminants in groundwater have varied since 2011 in any given monitoring well, as shown in **Table 6**. The groundwater contamination in Area 11 is primarily ETX compounds, with toluene being the most conspicuous because of the magnitude of its RG exceedances. Historically, several chlorinated VOCs have been detected over their RGs in samples collected from Area 11. A recurring issue has been that due to the high levels of ETX compounds, samples may require dilution, and this results in higher detection limits for other analytes, often above their RG. Due to this masking effect, it is not possible to determine if these compounds are actually present above their RGs in the diluted samples. At the end of 2020, monitoring well samples have been reliably analyzed for 1,4-dioxane for five events, one in 2019 and the four quarterly events in 2020. A discussion of its presence in these groundwater samples is included in this section.

An additional background well, MW-130A, was added to the monitoring well network in August 2020. This well is located immediately downgradient of Area 4. The compounds detected during the September and December sampling rounds were 1,1,1-trichloroethane (TCA), 1,1-dichloroethene (DCE) and 1,4-dioxane near the reporting limits, and well below the RGs.

Monitoring well MW-001 is also considered an upgradient, background well for Area 11. Samples collected from this well during five events from 2011 to 2012 contained TCA, DCE, and trichloroethene (TCE) at concentrations just above their respective RGs. Starting in December 2012, concentrations of these compounds decreased rapidly to low double- and single-digit levels that have remained consistently below RGs. It is believed that the decrease

in concentrations of chlorinated VOCs is attributable to the Area 4 hydraulic containment leachate component remedy that operated from December 2009 to October 2018. See **Figure 1** for the location of Area 4 relative to Area 11. The ETX compounds have been detected sporadically since 2011 at very low levels, near or below the analytical detection limits. The compound 1,4-dioxane was detected above the RG in two of the five sampling events and was not detected above the RG for the last two events of 2020. The average of the 1,4-dioxane detections for the five sampling events is approximately 8 micrograms per liter ( $\mu\text{g/L}$ ), which is slightly above the RG of  $7.7 \mu\text{g/L}$ .

Continuing in a generally hydrogeologic downgradient order, the well locations with the highest concentrations of contaminants continue to be MW-004A and MW-002. These wells are primarily contaminated with toluene, and concentrations have remained above the RG since 2011. In 2020, toluene concentrations in MW-002 for the first two sampling quarters decreased to approximately three times lower than the high concentrations in 2016 and 2017. The concentrations continued to decrease for the last two quarters to six times lower than previous high levels. The 2016 and 2017 high concentrations were measured prior to the Harrison Avenue construction and removal of impervious surfaces in the area.

Concentrations of ethylbenzene and xylenes in MW-002 remained at levels over their RGs since 2012, showing decreasing concentrations through 2019. Beginning in 2020, the concentrations of both compounds have gradually increased to levels two to three times the previous highest levels. In addition, benzene was detected once in a sample collected from MW-002 in June 2013, at a concentration just over its RG of  $5 \mu\text{g/L}$ . Vinyl chloride was detected in June 2020 at a concentration of  $4.4 \mu\text{g/L}$ , which is two times the RG of  $2.0 \mu\text{g/L}$ . The compound 1,4-dioxane was not detected above the RG of  $7.7 \mu\text{g/L}$  when analyzed in any of the 2019 or 2020 sampling events.

Toluene concentrations in samples collected from MW-004A through 2020 have decreased from a maximum of 230 times the RG ( $230,000 \mu\text{g/L}$ ) in June 2013, to the December 2020 concentration of 34 times the RG ( $34,200 \mu\text{g/L}$ ). The most extreme drop in concentration was between April 2016 and March 2017 when toluene concentrations dropped by half from 150 times the RG ( $150,000 \mu\text{g/L}$ ) in April 2016, to 79 times the RG ( $79,000$ ) in March 2017. Since 2017, toluene concentrations have fluctuated, but continued to decrease. Both ethylbenzene and xylenes concentrations have continued to decrease from their initial, higher levels in 2011 and 2012. The xylenes concentrations in samples collected from MW-004A have been below the RG of  $10,000 \mu\text{g/L}$  since 2012, and the ethylbenzene concentrations have been below the RG of  $700 \mu\text{g/L}$  since 2013. For the 2020 quarterly sampling events, the detection limits for 1,4-dioxane, ethylbenzene and xylenes were below their respective RGs, so the reporting of these compounds was not impacted by sample dilutions. The compound 1,4-dioxane was not detected above the RG in any of the 2019 or 2020 sampling events when analyzed.

Samples collected from MW-004B are similar to those collected from background wells MW-001 and MW-130A, with low double- and single-digit detections of several chlorinated compounds considered to be generally representative of site-wide groundwater quality. The steady decrease in TCA concentrations since 2011 is at least partially attributable to the Area 4 leachate component remedy. The compound 1,4-dioxane was detected above the RG in four of the five sampling events where analyzed. The average of the 1,4-dioxane detections for the five sampling events is approximately  $10 \mu\text{g/L}$ , which is above the RG of  $7.7 \mu\text{g/L}$  and is considered representative of 1,4-dioxane concentrations in site-wide groundwater.

MW-003, located just south of MW-002, has not shown significant ETX contamination since September 2012, however, the May 2019 and March 2020 sampling events showed a spike in total xylene concentration slightly over the RG. Additionally, in March 2020, the ethylbenzene concentration increased from well below the RG to twice the RG (1,500 µg/L). Ethylbenzene concentrations in this well are typically well below the RG of 700 µg/L. The average of the 1,4-dioxane detections for the five sampling events is approximately 8 µg/L, which is slightly above the RG of 7.7 µg/L. MW-003 is generally not believed to be directly downgradient of any known Area 11 source(s).

MW-007, the newest downgradient monitoring well installed in 2018, is directly west and downgradient of MW-002, as shown in **Figure 2**. To date, eight rounds of samples have been collected from this well with ethylbenzene above its RG for all sample events. Ethylbenzene concentrations have fluctuated ranging from 9.5 times the RG (6,700 µg/L) in November 2018, down to two times the RG (1,420 µg/L) a year later in November 2019 and increasing in 2020 to five times the RG (3,660 µg/L). Xylenes were detected above the RG for the first event and declined to levels approximately half the concentration of the RG during both 2019 sampling events. In 2020, xylenes increased and were detected at approximately two times the 2019 levels, but still below the RG. Toluene was detected below the RG in November 2018, but not detected in 2019 or 2020, despite being a relatively short distance downgradient from MW-002 with its high concentrations of toluene. Benzene was reported at eight times the RG in the field duplicate (44.3 µg/L) and was non-detect in the sample in December 2020. The results are considered estimated because the overall agreement between sample and field duplicate was poor. The data from 2018 and 2019 indicate that sample dilutions may have resulted in the masking of benzene detections in these earlier sampling events. The compound 1,4-dioxane was not detected above the RG when analyzed in any of the 2019 or 2020 sampling events.

Monitoring wells MW-005 and MW-006 are located downgradient of Area 11. Samples collected from these wells have contained a combination of site-wide chlorinated compounds, and ETX compounds at low double- and single-digit concentrations. The only compounds detected above RGs in either well are 1,4-dioxane, bromodichloromethane and benzene. Bromodichloromethane has been detected above the RG several times in both wells. Bromodichloromethane is a trihalomethane, generally referred to as a disinfection by-product resulting from chlorine treatment of drinking water that has been routinely detected above its RG in the Area 4 background monitoring well. Its detection is not considered to be attributable to either source area. Benzene has been detected once in samples collected from MW-005, and a number of times in MW-006, including once at a concentration above its RG of 5 µg/L in May 2019. Benzene is not known to be related to Area 11, but the possibility still exists that it is related. The average of the 1,4-dioxane detections for the five sampling events for MW-0005 is 6.8 µg/L and 4.5 µg/L for MW-0006, both below the RG of 7.7 µg/L.

### 3.3 Analytical Results for Attenuation Parameters

The analytical results for the attenuation parameters methane, nitrate, sulfate, and alkalinity are summarized in **Table 7**. The results of the dissolved ferrous iron field screening were provided earlier, in **Table 4**. The laboratory analytical results for the attenuation parameters were compared to the Groundwater Quality Standards for Class I: Potable Resource Groundwater, Title 35, 35 IAC, Section 620.410, where applicable. Attenuation parameter laboratory data was initially collected in August 2014, then sporadically until 2020, when data was collected for each quarterly sampling event.



The majority of wells had low- to non-detect levels of nitrate, with the exception of MW-005 and MW-001. In September and December 2020 nitrate slightly exceeded the RG in MW-001 and in December 2020 nitrate exceeded the RG in MW-005. Overall, alkalinity levels are consistent between wells over time, with MW-006 and MW-007 having slightly higher alkalinity concentrations in 2020 than previously detected. Sulfate levels are also fairly consistent between wells. Over time MW-002, MW-003 and MW-007 have had the lowest sulfate concentrations. No wells exceeded the RG for sulfate.

The majority of wells in Area 11 had very low-level methane detections, near or just above the reporting limits. The wells with relatively higher methane concentrations are MW-002, MW-003 and MW-007. MW-002 had detections ranging from approximately 3 mg/L in 2014, to 31 mg/L in December 2020. MW-003 had detections ranging from approximately 1 mg/L in 2014, to 12 mg/L in March 2020. The newest downgradient well, MW-007, had detections ranging from 4 mg/L to 31 mg/L during the period from November 2018 to December 2020.

The dissolved ferrous iron field screening data for the four sampling events in 2020 are presented in **Table 4**. The screening data results ranged from non-detect in MW-004B and MW-005, to 4.98 mg/L in MW-007. The wells that consistently showed dissolved ferrous iron concentrations above 2 mg/L during 2020 are MW-002, MW-003, MW-004A, MW-006 and MW-007.

In addition to changes resulting from the removal of the asphalt parking lot, the attenuation data provide support for degradation being a significant component of the decreasing concentrations observed at this site. MW-004A appears to be iron- to sulfate- reducing, while MW-002 and MW-003 are methanogenic. These conditions are indicative of significant degradation/consumption of carbon sources (e.g., in this case, the ETX contaminants) and depletion of electron donors, which provides further evidence of biodegradation.

## Section 4

### Conclusions

This report summarizes the information obtained during the quarterly groundwater monitoring events in 2020 at Area 11. Data are being collected to document the leachate component LTRA and to obtain data that will support further pre-design work on the soil component.

This was the first year of quarterly monitoring events planned to provide input to an updated CSM for Area 11. This update to the CSM was recommended due to the changes in surface features at Area 11, completed in 2016, and the completion of the remedial action at upgradient Area 4 in 2018. Based on groundwater data from 2017 to 2019, it was hypothesized that hydrogeologic and contaminant migration patterns have been influenced by these changes, altering the previous CSM. The 2020 data from this first year of quarterly monitoring support this hypothesis.

The data being collected will be used to evaluate the new CSM at the site. This evaluation will provide input to refine the remedial decisions for both the leachate and soil components of Area 11 and could eventually lead to a ROD modification. At this time, the data being collected is adequate for its intended use; however, additional pre-design investigation activities such as soil borings and monitoring well installation may be necessary to support a ROD modification.

#### 4.1 Hydraulic Results

Groundwater levels were measured for all four quarterly groundwater sampling events in 2020. The updated QAPP was approved in August 2020, which increased the number of wells and the coverage of the Area 11 monitoring well network for the third and fourth quarterly events. Potentiometric surface maps were prepared for these two quarterly events to provide a more accurate picture of the potentiometric surface. **Table 3** provides a summary of the groundwater elevation measurements for the 2020 sampling events, while incorporating the updated monitoring well survey information. **Figures 3 and 4** provide the potentiometric surface maps for the two 2020 quarterly events. Groundwater gradients across Area 11 are relatively flat, with a predominantly westerly flow direction. The gradients as the groundwater enters Area 11 on the eastern edge and trends northwest are slightly higher than the gradients as the groundwater exits the site after changing direction and trending west-southwest. The overall groundwater flow is consistent with historic trends.

#### 4.2 Monitoring Well Results

The Area 11 groundwater contamination is primarily located east of 11<sup>th</sup> street and in the upper portion of the aquifer. This is to be expected with the more prevalent, and lighter, ETX contaminants. This assessment has been confirmed by the previous pre-design investigations conducted in 2008 and 2013. The construction along Harrison and 11<sup>th</sup> was completed in 2016 which resulted in the removal of significant impervious surfaces in this area of higher contaminant concentrations. The most recent pre-design investigation in 2018 did not identify a vadose zone contaminant source in this area. An additional downgradient

monitoring well (MW-007) west of 11<sup>th</sup> Street was installed in October 2018, and an existing upgradient monitoring well (MW-130A), were both added to the monitoring well network.

After the installation and sampling of monitoring well MW-007, it was apparent that contamination above the RG is present in the groundwater downgradient of the source area. However, contaminant concentrations are attenuating rapidly with distance based on a comparison of concentrations in samples collected from MW-002 and MW-007 through the end of 2020.

In downgradient well MW-0007, ethylbenzene concentrations have fluctuated, ranging from 9.5 times the RG in November 2018, down to two times the RG in November 2019, and increasing in 2020 to five times the RG. Xylenes were detected above the RG in 2018, declined to below the RG in 2019, and increased in 2020 to approximately two times the 2019 levels, but still below the RG. Benzene was detected at eight times the RG in the field duplicate and was non-detect in the sample in December 2020. The overall agreement between sample and field duplicate was poor so the results remain suspect and are considered estimated. These fluctuations may be due to slugs of contamination moving slowly through the unconsolidated formation. This is likely due to the removal of the impervious surfaces, as well as the influence of precipitation events.

Additional evidence that the removal of the asphalt allows increased infiltration of precipitation, periodically flushing soil contaminants into the groundwater is seen at MW-003. This well, located just south of MW-002, has not shown significant ETX contamination since September 2012. However, the May 2019 and March 2020 sampling events showed a spike in total xylene concentration slightly over the RG. Additionally, in March 2020, the ethylbenzene concentration increased from well below the RG, to twice the RG, and decreased to below the RG for the last three sampling events.

From 2017 through 2020, toluene concentrations in samples collected from MW-004A and MW-002 have decreased significantly in a manner not previously observed at Area 11. It is believed that the cause of the sudden reduction in toluene concentrations in MW-004A between April 2016 and March 2017 and in MW-002 between November 2018 and May 2019, is related to the removal of all asphalt surfaces from the site in 2016. In contrast, concentrations of ethylbenzene and xylenes in MW-002 remained at levels over their RGs since 2012, showing decreasing concentrations through 2019. However, beginning in 2020, the concentrations of both compounds have gradually increased to levels two to three times the previous highest levels. This same pattern was observed at MW-007, as previously discussed. It is likely groundwater with much lower toluene concentrations that continues to migrate from MW-004A downgradient to MW-002. The groundwater patterns seen in MW-002 appear to migrate downgradient to MW-007. The exact attenuation and transport processes responsible for these significant changes in concentrations is not known but will continue to be evaluated.

The parameter 1,4-dioxane has been reliably analyzed for five sampling events in all wells, except MW-130A which has data for two events. The data from the 2020 quarterly sampling detected 1,4-dioxane at least once in all Area 11 monitoring network wells. Over the five events, five wells had concentrations that exceeded the RG at least once. The wells that have the most detections and highest concentrations are wells MW-001, MW-003 and MW-004B, which are all not directly impacted by the Area 11 source(s). Conversely, MW-002 and MW-



004A, which are the most impacted Area 11 wells, have the lowest 1,4-dioxane concentrations making it very unlikely that Area 11 is a source of 1,4-dioxane.

### 4.3 Recommendations

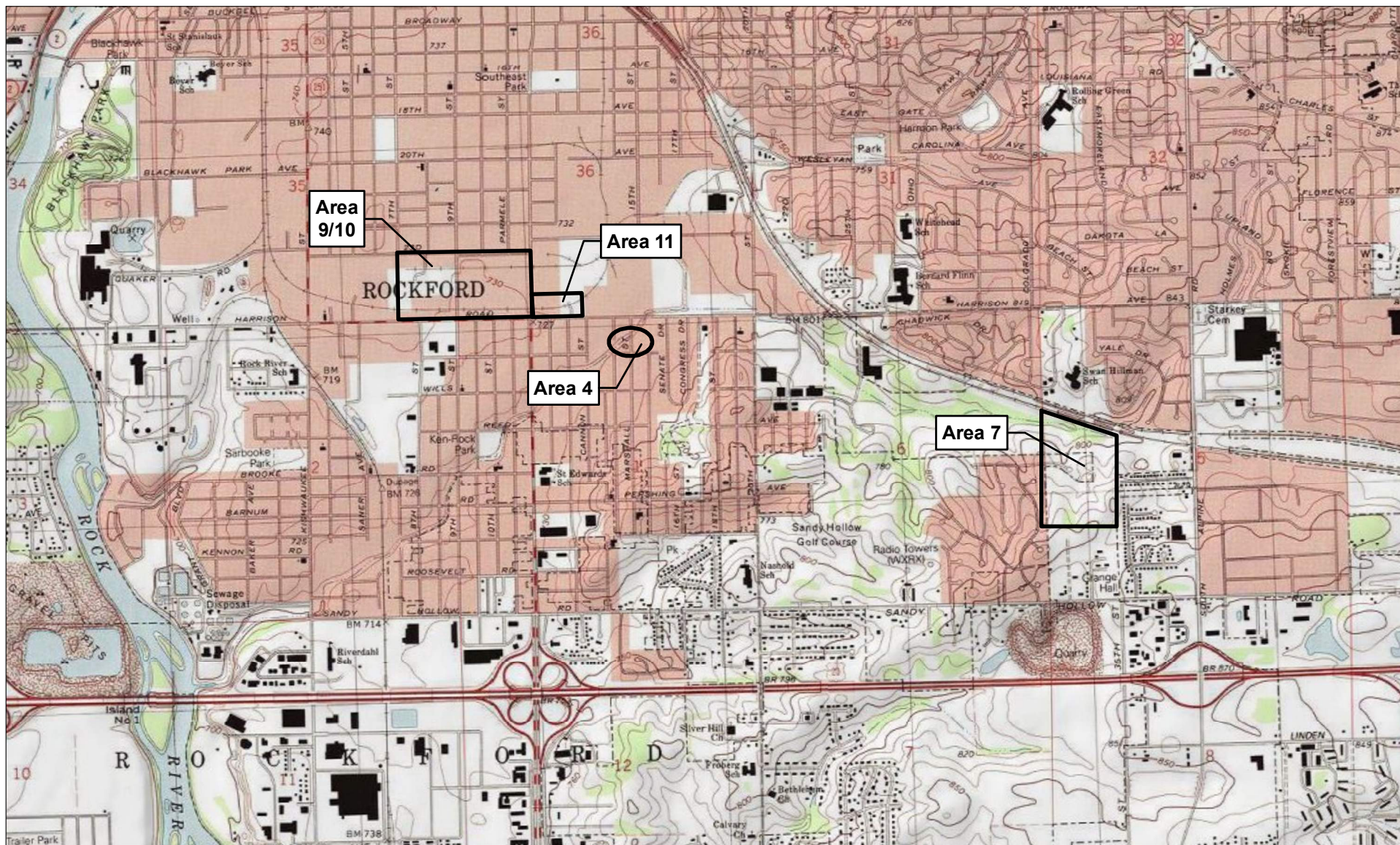
At this time, without the identification of a source, the plans for remedial design for the Area 11 soil component cannot be undertaken. However, due to the changes to the surficial conditions within Area 11, the overall conceptual site model has changed. This is evident from the recent significant changes in contaminant concentrations, as well as the trends in the monitoring wells that have historically shown the highest levels of site contaminants.

It is recommended that quarterly monitoring continue for at least another year. The VOCs and attenuation parameters should continue to be monitored, and the information gathered should be used to reevaluate the Area 11 conceptual site model and to assess the mechanisms that may be influencing the contaminant concentrations in groundwater within Area 11. This reevaluation will provide input to refine the remedial decisions for both leachate and soils within Area 11.


Based on the five monitoring events that included analysis of 1,4-dioxane, CDM Smith recommends that analysis of 1,4-dioxane be discontinued because results indicate Area 11 is not a source of 1,4-dioxane.

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# Figures



#### LEGEND

 Source Area

#### Service Layer Credits:

- Aerial Source: Esri, DigitalGlobe, GeoEye, i-cubed, USDA, USGS, AEX, Getmapping, Aerogrid, IGN, IGP, swisstopo, and the GIS User Community.
- Road Centerline Source: Census TIGER/Line Roads, 2020.

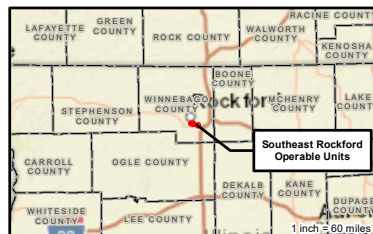
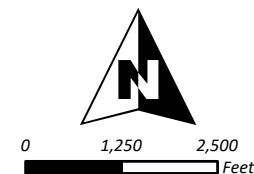


Figure 1 - Area Map







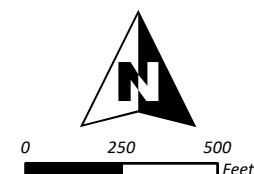
#### LEGEND

- Groundwater Sampling Location
- Water Level Measurement Location

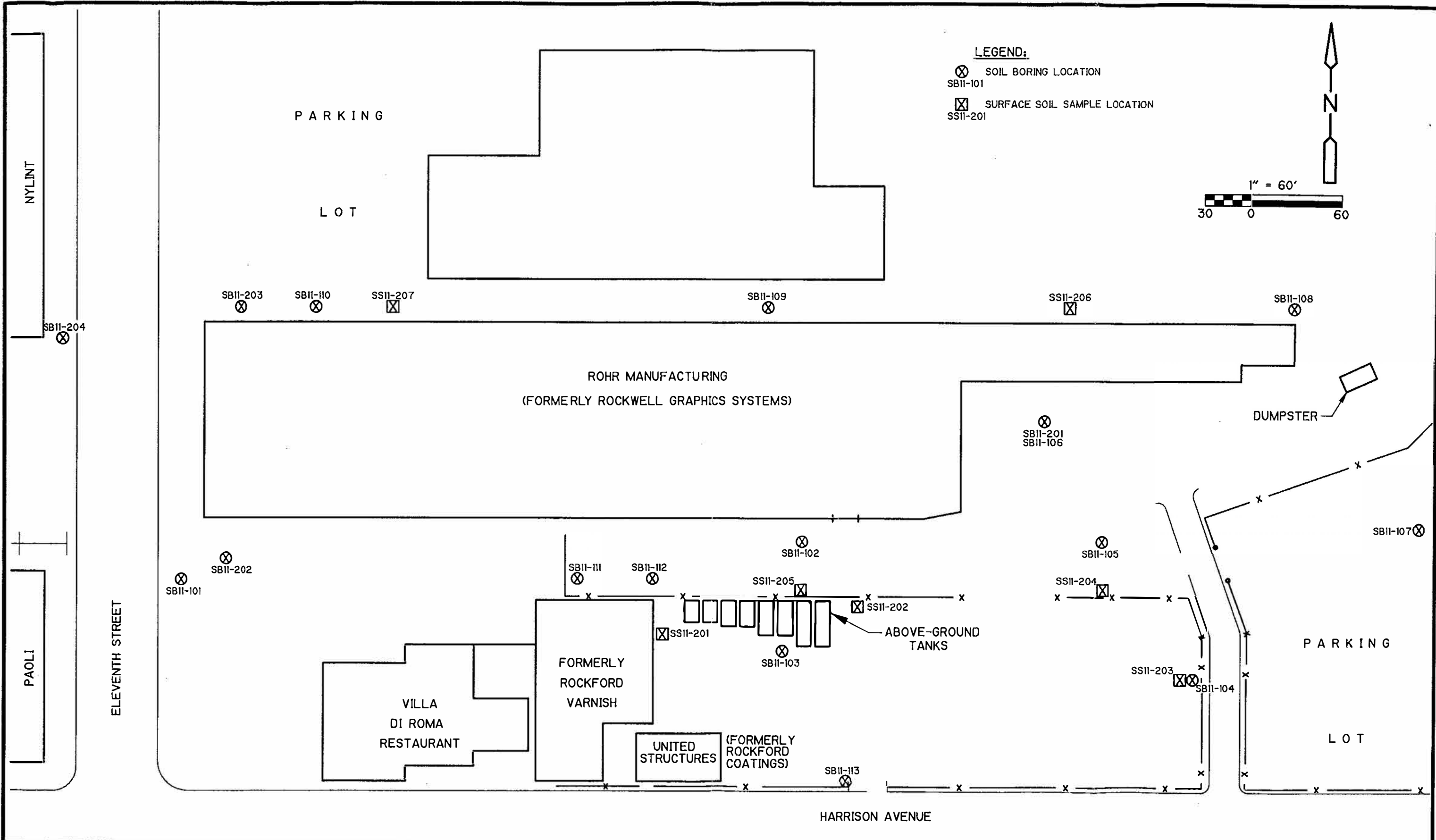
Figure 2 - Area 11 Monitoring Well Locations

#### Service Layer Credits:

- Aerial Source: Esri, DigitalGlobe, GeoEye, i-cubed, USDA, USGS, AEX, Getmapping, Aerogrid, IGN, IGP, swisstopo, and the GIS User Community.
- Road Centerline Source: Census TIGER/Line Roads, 2020.



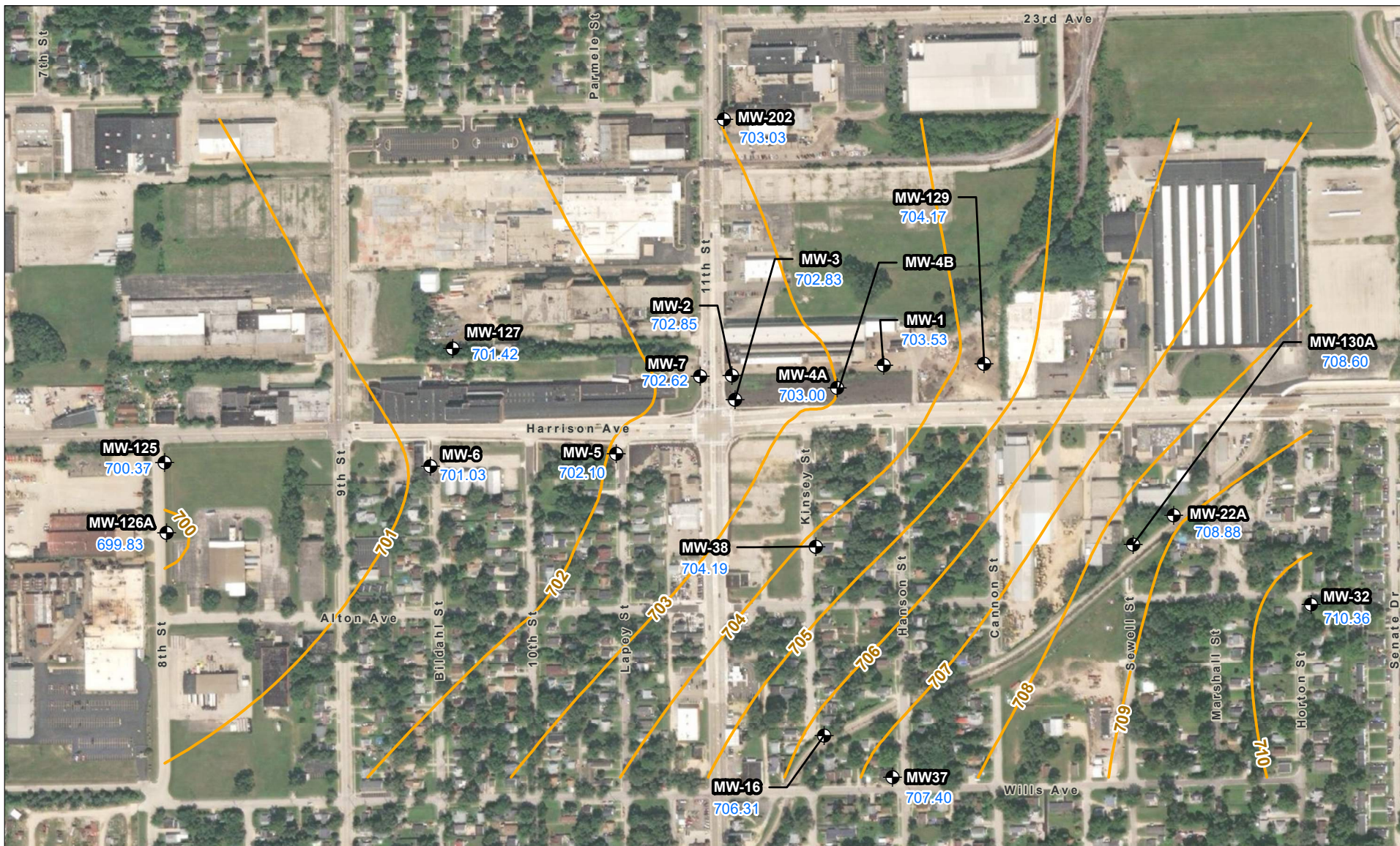
S:\1681\11110\RI\FINAL\AREA11\ 11\_sbloc 07/19/00 10:05:13 2:16:00 Selfgorb





SOUTHEAST ROCKFORD SOURCE CONTROL OPERABLE UNIT  
AREA 11 SOIL BORING AND SURFACE SOIL  
SAMPLE LOCATIONS

Figure No. 3-23





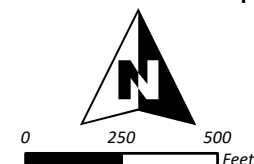
#### LEGEND

 Monitoring Well and Water Level Elevation  
 Groundwater Contour

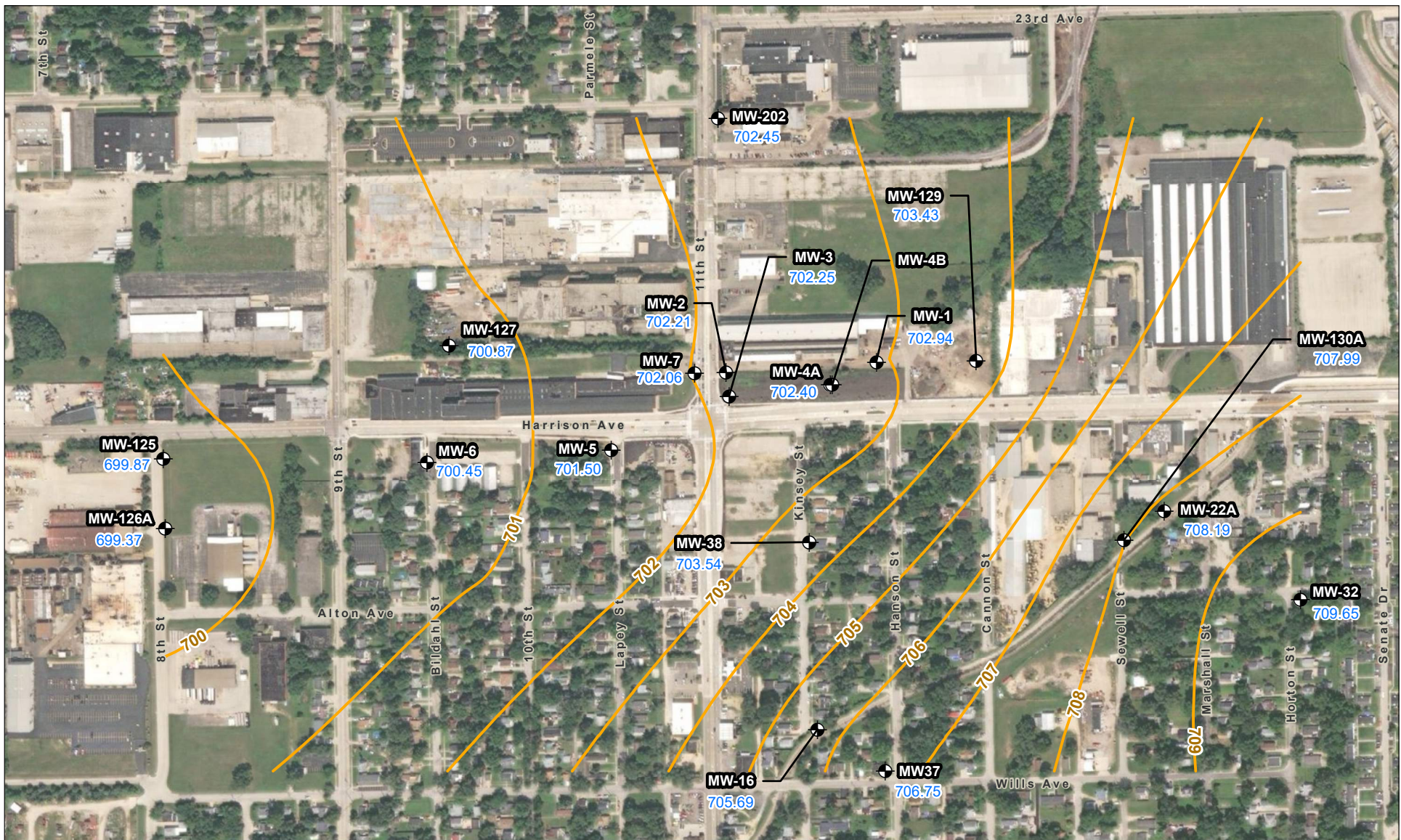
#### Service Layer Credits:

- Water level measurements taken October 15, 2020.
- Aerial Source: Esri, DigitalGlobe, GeoEye, i-cubed, USDA, USGS, AEX, Getmapping, Aerogrid, IGN, IGP, swisstopo, and the GIS User Community.
- Road Centerline Source: Census TIGER/Line Roads, 2020.

Figure 3 - Area 11 Third Quarter October 2020 Groundwater Potentiometric Surface Map







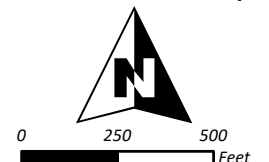
#### LEGEND

- Monitoring Well and Water Level Elevation  
 Groundwater Contour

#### Service Layer Credits:

- Water level measurements taken November 30, 2020.
- Aerial Source: Esri, DigitalGlobe, GeoEye, i-cubed, USDA, USGS, AEX, Getmapping, Aerogrid, IGN, IGP, swisstopo, and the GIS User Community.
- Road Centerline Source: Census TIGER/Line Roads, 2020.

Figure 4 - Area 11 Fourth Quarter November 2020 Groundwater Potentiometric Surface Map







# Tables

**Table 1**  
**2020 Groundwater Sampling Dates**  
**Source Area 11 2020 Groundwater Report**  
**Southeast Rockford Groundwater Contamination Superfund Site**

Sampling Event	MW-001	MW-002	MW-003	MW-004A	MW-004B	MW-005	MW-006	MW-007	MW-130A
1 <sup>st</sup> Quarterly 2020	3/3/2020	3/4/2020	3/4/2020	3/4/2020	3/3/2020	3/3/2020	3/3/2020	3/4/2020	NS
2 <sup>nd</sup> Quarterly 2020	6/9/2020	6/10/2020	6/10/2020	6/10/2020	6/9/2020	6/9/2020	6/9/2020	6/10/2020	NS
3 <sup>rd</sup> Quarterly 2020	9/9/2020	9/10/2020	9/10/2020	9/10/2020	9/9/2020	9/9/2020	9/9/2020	9/10/2020	9/9/2020
4 <sup>th</sup> Quarterly 2020	12/1/2020	12/2/2020	12/2/2020	12/2/2020	12/2/2020	12/1/2020	12/1/2020	12/2/2020	12/1/2020

NS – Not Sampled

**Table 2**  
**Source Area 11 Groundwater Monitoring Well Details**  
**Source Area 11 2020 Groundwater Report**  
**Southeast Rockford Groundwater Contamination Superfund Site**

Well Number	Depth to Screen Base from Ground Surface	Top of Casing Elevation	Top of Screen Elevation	Bottom of Screen Elevation	Screen Length	Aquifer Screened	Ground Surface Elevation
MW-130A	37.5	728.04	700.59	690.59	10	unconsolidated	728.09
MW-001	50	731.05	691.57	681.57	10	unconsolidated	731.44
MW-002	50	727.78	688.36	678.36	10	unconsolidated	728.18
MW-003	50	728.11	688.96	678.96	10	unconsolidated	728.55
MW-004A	40	729.66	700.12	690.12	10	unconsolidated	730.08
MW-004B	85	730.5	655.31	645.31	10	unconsolidated	730.39
MW-005	48	727.95	689.95	679.95	10	unconsolidated	728.35
MW-006	51	727.05	686.27	676.27	10	unconsolidated	727.41
MW-007	45	727.44	692.5	682.5	10	unconsolidated	727.8
MW-125*	46	727.75	691.9	681.9	10	unconsolidated	727.75
MW-126A*	55	727.84	682.9	672.9	10	unconsolidated	727.8
MW-127*	42	728.5	694.7	684.7	10	unconsolidated	726.54
MW-129*	32	731.6	705.1	700.1	5	unconsolidated	732.11
MW-202*	50	729.06	689.5	679.5	10	unconsolidated	729.19
MW-16*	53	725.33	677.8	672.8	5	unconsolidated	725.51
MW-22A*	38.5	730.35	702.2	692.2	10	unconsolidated	730.67
MW-32*	45	733.84	699.2	689.2	10	unconsolidated	734.16
MW-37*	44	725.05	686.1	681.1	5	unconsolidated	725.08
MW-38*	48	728.28	685.2	680.2	5	unconsolidated	728.79

**Table 3**  
**2020 Observed Groundwater Elevations**  
**Source Area 11 2020 Groundwater Report**  
**Southeast Rockford Groundwater Contamination Superfund Site**

Well ID	Top of Casing Elevation (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevation (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevation (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevation (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevation (ft AMSL)
		March 2, 2020		June 8, 2020		October 15, 2020		November 30, 2020	
MW-001	731.05	26.92	704.13	26.36	704.70	27.52	703.53	28.11	702.94
MW-002	727.78	24.39	703.39	23.78	704.00	24.93	702.85	25.57	702.21
MW-003	728.11	24.68	703.43	24.06	704.05	25.28	702.83	25.86	702.25
MW-400A	729.66	26.09	703.57	25.46	704.20	26.66	703.00	27.26	702.40
MW-004B*	730.50	25.68	704.82	25.10	705.40	26.27	704.23	26.88	703.62
MW-005	727.95	25.33	702.62	24.64	703.31	25.85	702.10	26.45	701.50
MW-006	727.05	25.49	701.56	24.76	702.29	26.02	701.03	26.60	700.45
MW-007	727.44	24.19	703.25	23.59	703.85	24.82	702.62	25.38	702.06
MW-130A	728.04	NA	NA	NA	NA	19.44	708.60	20.05	707.99
MW-16	725.33	18.60	706.73	17.92	707.41	19.02	706.31	19.64	705.69
MW-127	728.50	26.54	701.96	25.82	702.68	27.08	701.42	27.63	700.87
MW-129	731.60	26.95	704.65	26.39	705.21	27.43	704.17	28.17	703.43
MW-202	729.06	23.39	705.67	24.81	704.25	26.03	703.03	26.61	702.45
MW-32	733.84	NA	NA	NA	NA	23.48	710.36	24.19	709.65
MW-22A	730.35	NA	NA	NA	NA	21.47	708.88	22.16	708.19
MW-37	725.05	NA	NA	22.96	702.09	17.65	707.40	18.30	706.75
MW-38	728.28	NA	NA	29.85	698.43	24.09	704.19	24.74	703.54
MW-125	727.75	NA	NA	NA	NA	27.38	700.37	27.88	699.87
MW-126A	727.84	NA	NA	NA	NA	28.01	699.83	28.47	699.37

**Notes:**

\* well not included in potentiometric surface maps

AMSL = above mean sea level

BTOC = below top of casing

ft = feet

NA = not available

**Table 4**  
**2020 Final Stabilized Field Parameter Readings for Monitoring Well Purging**  
**Source Area 11 2020 Groundwater Report**  
**Southeast Rockford Groundwater Contamination Superfund Site**

	Flowrate mL/min	pH	Specific Cond. (mS/cm)	Turbidity (NTU)	Dissolved Oxygen (mg/L)	Temp (°C)	ORP (mV)	Dissolved Ferrous Iron (mg/L)	Purged Min.
<b>Mar-20</b>									
MW-001	NA	7.27	1.41	9.3	5.06	12.08	118	0.04	50
MW-002	NA	6.54	1.36	22	0.83	13.61	-116	1.76	35
MW-003	NA	6.62	1.35	17.7	3.39	12.12	-122	1.32	30
MW-004A	NA	6.74	1.09	5.7	3.53	12.44	-134	2.4	25
MW-004B	NA	7.42	1.18	100	1.66	12.26	37	0.1	50
MW-005	NA	7.47	1.26	20.2	4.6	10.36	70	0.002	55
MW-006	NA	7.38	1.35	0.6	0.78	12.92	-110	1.72	25
MW-007	NA	6.75	0.868	10.1	6.75	10.3	-131	1.8	35
<b>Jun-20</b>									
MW-001	NA	6.75	1.21	58	8.48	17.73	71	0.1	50
MW-002	250	6.65	1.33	19	2.5	17.99	-106	4.58	35
MW-003	400	6.94	1.26	11.1	3.54	15.26	-131	3.48	50
MW-004A	400	7.01	1.12	2.6	5.41	14.22	-272	2.39	50
MW-004B	500	6.78	1.18	26.4	3.11	17.31	5	0.18	35
MW-005	450	6.8	1.43	32.8	5.24	17.71	134	ND	25
MW-006	400	6.8	1.49	4.8	3.11	15.71	-115	3.84	35
MW-007	350	6.94	0.992	4	4.2	16.42	-140	2.39	25
<b>Sep-20</b>									
MW-001	300	6.93	1.28	81	3.84	14.31	64	0.02	45
MW-002	250	6.71	1.37	13.3	0.33	15.73	-116	2.14	40
MW-003	400	7.03	1.26	9.9	0.44	14.13	-140	2.52	45
MW-004A	325	7.13	1.29	9.8	4.5	14.44	-226	1.12	60
MW-004B	300	7.04	1.25	13.1	2.8	14.05	41	ND	55
MW-005	425	7	1.26	10	3.73	14.43	78	ND	50
MW-006	275	6.94	1.36	12.3	0.49	14.88	-133	2.19	30
MW-007	400	6.57	1.94	9.9	0.34	14.82	-126	4.98	45
MW-130A	425	6.69	1.18	9.8	2.44	13.85	24	NA	45
<b>Dec-20</b>									
MW-001	475	8.02	1.19	12	5.69	12.08	20	ND	90
MW-002	350	7.21	1.43	199	0.39	14.18	-109	2.7	60
MW-003	375	7.55	1.34	7	0.62	11.75	-158	3.85	45
MW-004A	330	7.43	1.34	3.6	0.54	12.34	-276	2.61	40
MW-004B	400	7.39	1.4	32	4.15	10.93	94	0.73	65
MW-005	500	7.16	1.49	39.9	4.85	12.84	160	0.03	60
MW-006	500	7.92	1.38	2.7	1.33	12.66	-150	2.96	45
MW-007	450	7.02	1.38	20.7	0.5	12.57	-128	2.86	40
MW-130A	350	7.9	1.17	27.2	2.19	10.71	-21	0.19	60

NA = Not Available

ND = Non-Detect

Table 5  
VOC Compounds Detected 2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location Sample ID Sample Date		A11-MW001 A11-MW001-200303 3/3/2020	A11-MW001 A11-MW001-200609 6/9/2020	A11-MW001 A11-MW001-200909 9/9/2020	A11-MW001 A11-MW001-201201 12/1/2020
Analyte Name	RG				
1,1,1-Trichloroethane	200	6.74 J	8.9	7.58 J	9.02
1,1-Dichloroethane	1400	4.51	7.5	5.16	4.94
1,1-Dichloroethene	7	2.00 U	1.4	2.00 U	2.00 U
1,4-Dioxane	7.7	6.85	14.1	0.205 U	5
cis-1,2-Dichloroethene	70	2.00 U	1.4	2.00 U	2.00 U
Tetrachloroethene	5	2.00 U	1	2.00 U	2.00 U
trans-1,2-Dichloroethene	100	2.00 U	0.17 J	2.00 U	2.00 U
Trichloroethene	5	2.00 U	2.5	2.41	2.15

Station Location Sample ID Sample Date		A11-MW002 A11-MW002-200304 3/4/2020	A11-MW002 A11-MW002-200610 6/10/2020	A11-MW002 A11-MW002-200910 9/10/2020	A11-MW002 A11-MW002-201201 12/1/2020
Analyte Name	RG				
1,1,1-Trichloroethane	200	100 U	1 J	50.0 U	50.0 U
1,1-Dichloroethane	1400	100 U	6.4	50.0 U	50.0 U
1,1-Dichloroethene	7	100 U	5 U	50.0 U	50.0 U
1,2,4-Trimethylbenzene	--	822	NA	622	588
1,2-Dichlorobenzene	600	100 U	5	50.0 U	50.0 U
1,3,5-Trimethylbenzene	--	285	NA	202	161
1,4-Dioxane	7.7	3.31	4.03	2.9	1.1
Acetone	6300	625 UJ	2.7 J	312 U	312 U
cis-1,2-Dichloroethene	70	100 U	32	50.0 U	50.0 U
Cyclohexane	--	NA	120	NA	NA
Ethyl Benzene	700	6840	6400	8260	10200
Isopropyl Benzene	700	121	98	90	78.1
Methyl Acetate	--	NA	11	NA	NA
Methylcyclohexane	--	NA	570 J	NA	NA
Naphthalene	140	100 U	NA	55.2	58.5
n-Propylbenzene	--	215	NA	129	87.4
Toluene	1000	78600	68000 J	39300	33200
Trichloroethene	5	100 U	0.86 J	50.0 U	50.0 U
Vinyl Chloride	2	100 U	4.4 J	50.0 U	50.0 U
Xylenes - Total	10000	24900	24900	32820	38040

Notes:  
All results in microgram per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded result exceeds remediation goal  
U = Not detected above the reported limit  
J = Estimated result  
N = Normal Sample  
\*\* = Field Duplicate Sample



Table 5  
VOC Compounds Detected 2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location Sample ID Sample Date		A11-MW003 A11-MW003-200304 3/4/2020	A11-MW003 A11-MW003-200610 6/10/2020	A11-MW003 A11-MW003-200910 9/10/2020	A11-MW003 A11-MW003-201201 12/1/2020
Analyte Name	RG				
1,1,1-Trichloroethane	200	20.0 U	1.3 J	10.0 U	10.0 U
1,1-Dichloroethane	1400	20.0 U	6.9	10.0 U	10.0 U
1,2,4-Trimethylbenzene	--	329	NA	113	178
1,3,5-Trimethylbenzene	--	121	NA	34.8	55
1,4-Dioxane	7.7	8.57	9.58	7.23	4.6
4-Methyl 2-Pentanone	--	50.0 U	2.2 J	25.0 U	25.0 U
Chloroethane	--	20.0 U	1.9 J	10.0 U	10.0 U
Cyclohexane	--	NA	7.8	NA	NA
Ethyl Benzene	700	1500	430 J	201	256
Isopropyl Benzene	700	74.4	45	28.3	38.5
Methylcyclohexane	--	NA	59	NA	
Naphthalene	140	29.1	NA	10.0 U	16.6
n-Butylbenzene	--	23.7	NA	10.0 U	13.1
n-Propylbenzene	--	76	NA	28.2	37.3
sec-Butylbenzene	--	27.8	NA	12	15
Toluene	1000	38.4	7.6 J	10.0 U	10.0 U
Xylenes - Total	10000	13000	5105.2	2430	6310

Station Location Sample ID Sample Date		A11-MW004A A11-MW004A-200304 3/4/2020	A11-MW004A A11-MW004A-200610 6/10/2020	A11-MW004A A11-MW004A-200910 9/10/2020	A11-MW004A A11-MW004A-201201 12/2/2020
Analyte Name	RG				
1,1,1-Trichloroethane	200	50.0 U	3.7 J	50.0 U	50.0 U
1,4-Dioxane	7.7	1.41	1.51	1.09	0.15 J
Cyclohexane	--	NA	0.83 J	NA	NA
Ethyl Benzene	700	260	330 J	365	331
Isopropyl Benzene	700	50.0 U	2.1 J	50.0 U	50.0 U
Methylcyclohexane	--	NA	21	NA	NA
Tetrachloroethene	5	50.0 U	5.3	50.0 U	50.0 U
Toluene	1000	45300	52000	42600	34200
Trichloroethene	5	50.0 U	1 J	50.0 U	50.0 U
Xylenes - Total	10000	414	531 J	604.6	541.2

Notes:  
All results in microgram per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded result exceeds remediation goal  
U = Not detected above the reported limit  
J = Estimated result  
N = Normal Sample  
\*\* = Field Duplicate Sample





Table 5  
VOC Compounds Detected 2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location Sample ID Sample Date		A11-MW004B A11-MW004B-200303 3/3/2020	A11-MW004B A11-MW004B-200609 6/9/2020	A11-MW004B A11-MW004B-200909 9/9/2020	A11-MW004B A11-MW004B-201201 12/1/2020
Analyte Name	RG				
1,1,1-Trichloroethane	200	5.29 J	5.5	4.93 J	5.61
1,1-Dichloroethane	1400	5.86	6.3	5.34	5.67
1,1-Dichloroethene	7	2.00 U	0.95	2.00 U	2.00 U
1,4-Dioxane	7.7	9.75	11.7	7.86	6.3
cis-1,2-Dichloroethene	70	2.00 U	1.3	2.00 U	2.00 U
Tetrachloroethene	5	2.00 U	0.4 J	2.00 U	2.00 U
Toluene	1000	2.00 U	1.6	2.00 UJ	2.00 U
trans-1,2-Dichloroethene	100	2.00 U	0.16 J	2.00 U	2.00 U
Trichloroethene	5	2.00 U	1.4	2.00 U	2.00 U

Station Location Sample ID Sample Date		A11-MW005 A11-MW005-200303 3/3/2020	A11-MW005 A11-MW005-200609 6/9/2020	A11-MW005 A11-MW005-200909 9/9/2020	A11-MW005 A11-MW005-201201 12/1/2020
Analyte Name	RG				
1,1,1-Trichloroethane	200	2.92 J	4.5	5.56 J	4.9
1,1-Dichloroethane	1400	3.77	6.4	9.11	7.01
1,1-Dichloroethene	7	2.00 U	1.1	2.00 UJ	2.00 U
1,4-Dioxane	7.7	5.35	8.83	8.18 J	4.5 J
Bromodichloromethane	0.2*	2.00 U	0.4 J	2.00 U	2.00 U
Chloroform	70	2.00 U	0.45 J	2.00 U	2.00 U
cis-1,2-Dichloroethene	70	2.00 U	1.3	2.00 U	2.00 U
Dibromochloromethane	140*	2.00 U	0.18 J	2.00 U	2.00 U
Tetrachloroethene	5	2.00 U	0.39 J	2.00 UJ	2.00 U
trans-1,2-Dichloroethene	100	2.00 U	0.15 J	2.00 UJ	2.00 U
Trichloroethene	5	2.00 U	0.89	2.00 UJ	2.00 U

Station Location Sample ID Sample Date		A11-MW006 A11-MW006-200303 3/3/2020	A11-MW006 A11-MW006-200609 6/9/2020	A11-MW006 A11-MW006-200909 9/9/2020	A11-MW006 A11-MW006-201201 12/1/2020
Analyte Name	RG				
1,1-Dichloroethane	1400	2.00 U	0.11 J	2.00 U	2.00 U
1,4-Dioxane	7.7	1.54	7.53	8.42	4.1
Benzene	5	2.62	2	2.28	2.82
Chloroethane	--	2.00 U	0.44 J	2.00 U	2.00 U
Cyclohexane	--	NA	0.35 J	NA	NA
Isopropyl Benzene	700	2.00 U	0.14 J	2.00 U	2.00 U
Trichloroethene	5	2.00 U	0.14 J	2.00 U	2.00 U

Notes:  
All results in microgram per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded result exceeds remediation goal  
U = Not detected above the reported limit  
J = Estimated result  
N = Normal Sample  
\*\* = Field Duplicate Sample



Table 5  
VOC Compounds Detected 2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location Sample ID Sample Date		A11-MW007 A11-MW007-200304 3/4/2020	A11-MW007** A11-MW007-200304-D 3/4/2020	A11-MW007 A11-MW007-200610 6/10/2020	A11-MW007** A11-MW007-200610-D 6/10/2020	A11-MW007 A11-MW007-200910 9/10/2020	A11-MW007** A11-MW007-200910-D 9/10/2020	A11-MW007 A11-MW007-201201 12/2/2020	A11-MW007** A11-MW007-201201-D 12/2/2020
Analyte Name	RG								
1,2,4-Trimethylbenzene	--	22.6	21	NA	NA	53.5	55.7	131	169
1,3,5-Trimethylbenzene	--	4.84	4.57	NA	NA	11.1	11.6	14.4 J	56.7 J
1,4-Dioxane	7.7	3.38	3.3	0.205 U	0.205 U	0.212 U	0.203 U	0.069 J	0.19 U
Benzene	5	4.00 U	4.00 U	5 U	5 U	10.0 U	10.0 U	10.0 UJ	44.3 J
Ethyl Benzene	700	959	863	820	810	2630	2680	3300	3660
Isopropyl Benzene	700	12	11	6.5	6.5	86.1	89.1	109 J	486 J
Methylcyclohexane	--	NA	NA	2.7 J	2.6 J	NA	NA	NA	NA
Naphthalene	140	4.00 U	4.00 U	NA	NA	11.4	13.2	34.0 J	97.3 J
n-Butylbenzene	--	4.00 U	4.00 U	NA	NA	11.3	12.4	19.9 J	66.9 J
n-Propylbenzene	--	6.68	6.03	NA	NA	82.4	84.7	104 J	454 J
sec-Butylbenzene	--	4.00 U	4.00 U	NA	NA	10.8	11.3	17.5 J	68.4 J
Tetrachloroethene	5	4.00 U	4.00 U	1 J	0.89 J	10.0 U	10.0 U	10.0 U	10.0 U
Xylenes - Total	10000	3050	2800	2600	2600	7600	7920	7390	8100

Station Location Sample ID Sample Date		A11-MW130A A11-MW130A-200909 9/9/2020	A11-MW130A A11-MW130A-201201 12/1/2020
Analyte Name	RG		
1,1,1-Trichloroethane	200	3.51 J	3.51
1,1-Dichloroethane	1400	4.11	3.77
1,4-Dioxane	7.7	6.1	4

Notes:  
All results in microgram per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded result exceeds remediation goal  
U = Not detected above the reported limit  
J = Estimated result  
N = Normal Sample  
\*\* = Field Duplicate Sample



Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW001 E52H2 A11-MW001-110419 4/19/2011	A11-MW001 E52L5 A11-MW001-110720 7/20/2011	A11-MW001 E52R7 A11-MW001-120111 1/11/2012	A11-MW001 E52S7 A11-MW001-04/02/2012 4/2/2012	A11-MW001 E3XB9 A11-MW001-120918 9/18/2012	A11-MW001 E3XC9 A11-MW001-121204 12/4/2012	A11-MW001 E3XF8 A11-MW001-130314 3/14/2013	A11-MW001 E3XG8 A11-MW001-130625 6/25/2013	A11-MW001 E3XP2 A11-MW001-140806 8/6/2014	A11-MW001 E3XQ3 A11-MW001-141217 12/17/2014	A11-MW001 E3XX1 A11-MW001-150520 5/20/2015
Analyte Name	RG											
1,1,1-Trichloroethane	200	240	210 D	200	210 D	150 J	34	36 D	47 D	15	18	17 J-
1,1,2-Trichloroethane	5	20 U	0.41 J	5 U	0.44 J	5 U	5 U	0.5 U	0.068 J	0.5 U	0.5 UJ	0.5 UJ
1,1-Dichloroethane	1400	24	21	25	18	7.6	2.9 J	4.3	5.7	3	4.1 J	5 J-
1,1-Dichloroethene	7	11 J	11	5 U	11	9.4	5 U	3.2	3.8	1.1 J	2.1	1.9 J-
1,4-Dioxane	7.7	400 R	100 R	100 R	100 R	100 R	100 U	NA	NA	NA	NA	NA
Carbon Tetrachloride	5	20 U	5 U	5 U	5.0 U	5 U	3.8 J	0.5 U	0.5 U	0.5 U	0.5 U	2.3 J-
cis-1,2-Dichloroethene	70	9.4 J	7.9	9.2	5.8	2.9 J	1.4 J	2.3	3	1.5 J	1.6	1.8 J-
Ethyl Benzene	700	20 U	0.25 J	5 U	5.0 U	5 U	5 U	0.14 J	0.12 J	0.18 J	0.5 UJ	0.5 UJ
Isopropyl Benzene	700	20 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.26 J	0.5 U	0.5 UJ	0.5 UJ
Tetrachloroethene	5	3.5 J	3.7 J	4.6 J	4.0 J	4.4 J	2.7 J	1.8	2.8	1.1	1.3 J	1.3 J-
Toluene	1000	20 U	5 U	5 U	0.68 J	0.92 J	5 U	0.5 U	0.52	1 U	0.5 UJ	0.5 UJ
trans-1,2-Dichloroethene	100	20 U	5 U	5 U	5.0 U	5 U	5 U	0.1 J	0.13 J	0.5 U	0.5 U	0.13 J-
Trichloroethene	5	8.6 J	6.1	4.7 J	4.2 J	4.1 J	2.3 J	2.2	4	1.5	1.7 J	2.1 J-
Trichlorofluoromethane (Freon 11)	2100	20 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.19 J-
Vinyl Chloride	2	20 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 UJ
Xylenes (Total)	10000	40 U	0.5 J	5 U	5 U	5 U	5 U	0.47 J	0.65 J	0.26 J	0.5 UJ	0.5 UJ

Station Location EPA Sample ID Sample ID Sample Date		A11-MW001 E3XZ8 A11-MW001-160406 4/6/2016	A11-MW001 E3Y44 A11-MW001-170309 3/9/2017	A11-MW001** E3Y45 A11-MW001-170309-D 3/9/2017	A11-MW001 E3YA2 A11-MW001-181113 11/13/2018	A11-MW001 E3YF1 A11-MW001-190520 5/20/2019	A11-MW001 A11-MW001-191112 11/12/2019	A11-MW001 A11-MW001-200303 A11-MW001-200303 3/3/2020	A11-MW001 E3YG2 A11-MW001-200609 6/9/2020	A11-MW001 A11-MW001-200909 A11-MW001-200909 9/9/2020	A11-MW001 E3YJ1 A11-MW001-201201 12/1/2020
Analyte Name	RG										
1,1,1-Trichloroethane	200	17	11	12	9.3	22	7.21	6.74 J	8.9	7.58 J	9.02
1,1,2-Trichloroethane	5	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
1,1-Dichloroethane	1400	6.5	6.2	6.9	9.8	20	5.25	4.51	7.5	5.16	4.94
1,1-Dichloroethene	7	0.5 U	1.4	1.5	1.7	4.6	2.00 U	2.00 U	1.4	2.00 U	2.00 U
1,4-Dioxane	7.7	NA	NA	NA	NA	NA	12.2	6.85	14.1	0.205 U	5
Carbon Tetrachloride	5	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
cis-1,2-Dichloroethene	70	1.4	1.3	1.4	1.7	3.4	2.00 U	2.00 U	1.4	2.00 U	2.00 U
Ethyl Benzene	700	0.15 J	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Isopropyl Benzene	700	0.5 U	0.5 U	0.5 U	0.5 U	0.5 R	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Tetrachloroethene	5	1.2	0.81	0.89	1	2.7	2.00 U	2.00 U	1	2.00 U	2.00 U
Toluene	1000	0.5 U	0.09 J	0.1 J	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 UJ	2.00 U
trans-1,2-Dichloroethene	100	0.17 J	0.14 J	0.16 J	0.25 J	0.46 J	2.00 U	2.00 U	0.17 J	2.00 U	2.00 U
Trichloroethene	5	2.8	2.2	2.4	3.7	4.7	2.00 U	2.00 U	2.5	2.41	2.15
Trichlorofluoromethane (Freon 11)	2100	0.32 J	0.5 U	0.5 U	0.15 J	0.32 J	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Vinyl Chloride	2	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Xylenes (Total)	10000	0.85	0.5 U	0.11 J	0.5 U	0.3 J+	4.00 U	4.00 U	0.5 U	4.00 U	4.00 U

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
U = Not detected at value shown  
J = Estimated result      J- = Estimated result biased low      J+ = Estimated result biased high  
R = Rejected



Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW002 E52K7 A11-MW002-110420 4/20/2011	A11-MW002 E52N5 A11-MW002-110721 7/21/2011	A11-MW002 E52S4 A11-MW002-120112 1/12/2012	A11-MW002 E52S8 A11-MW002-04/03/2012 4/3/2012	A11-MW002 E3XC0 A11-MW002-120919 9/19/2012	A11-MW002 E3XD0 A11-MW002-121205 12/5/2012	A11-MW002 E3XF9 A11-MW002-130314 3/14/2013	A11-MW002 E3XH0 A11-MW002-130626 6/26/2013	A11-MW002 E3XP7 A11-MW002-140807 8/7/2014	A11-MW002 E3XQ7 A11-MW002-141217 12/17/2014	A11-MW002 E3XX5 A11-MW002-150521 5/21/2015
Analyte Name	RG											
1,1,1-Trichloroethane	200	5 U	50 U	500 U	250 U	1000 U	4000 U	110 J	34	6300 U	1000 U	73 J-
1,1,2-Trichloroethane	5	10	50 U	500 U	250 U	1000 U	4000 U	5 U	5 U	6300 U	1000 U	1000 UJ
1,1-Dichloroethane	1400	9.5	50 U	500 U	78 J	1000 U	4000 U	76	46	6300 U	1000 U	1000 UJ
1,1-Dichloroethene	7	5 U	50 U	500 U	250 U	1000 U	4000 U	43	5 U	6300 U	1000 U	1000 UJ
1,2,4-Trimethylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,2-Dichlorobenzene	600	5 U	50 U	500 U	250 U	1000 U	4000 U	5 U	5 U	6300 U	1000 U	1000 UJ
1,2-Dichloropropane	5	7.6	50 U	500 U	250 U	1000 U	4000 U	5 UJ	5 UJ	6300 U	1000 U	1000 UJ
1,3,5-Trimethylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,4-Dioxane	7.7	100 R	1000 R	10000 R	5000 R	20000 R	80000 U	100 R	100 R	130000 R	20000 R	20000 R
2-Butanone	4200	10 U	100 U	1000 U	500 U	2000 U	8000 U	58	27	13000 U	2000 U	2000 UJ
2-Hexanone	--	10 U	100 U	1000 U	500 U	2000 U	8000 U	10 U	10 U	13000 U	2000 U	4900 J-D
4-Methyl 2-Pentanone	--	10 U	100 U	1000 U	500 U	2000 U	8000 U	16	2.2 J	13000 U	2000 U	2000 UJ
Acetone	6300	20 U	100 U	1000 U	1000 U	2000 U	8000 U	10 U	7.1 J	13000 U	2000 U	2000 UJ
Benzene	5	5 U	50 U	500 U	250 U	1000 U	4000 U	5 UJ	5.3	6300 U	1000 U	1000 UJ
Chloroethane	--	11	50 U	500 U	250 U	1000 U	4000 U	5 U	5 U	6300 UJ	1000 U	1000 UJ
cis-1,2-Dichloroethene	70	5 U	50 U	500 U	100 J	1000 U	4000 U	160	69	6300 U	1000 U	88 J-
Cyclohexane	--	98	81	500 U	82 J	1000 U	4000 U	100 J	170 J	6300 U	1000 U	1000 UJ
Dichlorodifluoromethane (Freon 12)	1400	5.7	50 U	500 U	250 U	1000 U	4000 U	5 U	5 U	6300 UJ	1000 U	1000 UJ
Ethyl Benzene	700	2700 JD	2000 D	3700	2700	1500	3900 J	1400 J	3400 D	3700 J	3100	7200 J-D
Isopropyl Benzene	700	75	94	500 U	72 J	1000 U	4000 U	53 J	85 J	6300 U	1000 U	77 J-
Methyl Acetate	--	5 U	50 U	500 U	250 U	1000 U	4000 U	5 U	2.8 J	6300 U	1000 U	1000 UJ
Methylcyclohexane	--	71 JD	420	280 J	340	230 J	4000 U	440 J	590 D	6300 U	1000 U	470 J-
Naphthalene	140	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
n-Propylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Styrene	100	5 U	50 U	500 U	250 U	1000 U	4000 U	5 R	5 UJ	6300 U	1000 U	1000 UJ
Tetrachloroethene	5	5 U	50 U	500 U	250 U	1000 U	4000 U	5 R	2.4 J	6300 U	1000 U	1000 UJ
Toluene	1000	110	50 U	NA	160000 D	360 J	220000 D	200000 D	210000 D	150000	190000 D	110000 J-D
Trichloroethene	5	5 U	50 U	500 U	250 U	1000 U	4000 U	5 R	7.6 J	6300 U	1000 U	1000 UJ
Vinyl Chloride	2	5 U	50 U	500 U	250 U	1000 U	4000 U	12	5 U	6300 U	1000 U	1000 UJ
Xylenes (Total)	10000	9010 JD	8371 D	14700	11500	11000	16400	13100	14400	15100	12700	26300

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
U = Not detected at value shown  
J = Estimated result      J- = Estimated result biased low      J+ = Estimated result biased high  
R = Rejected



Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW002 E3XZ9 A11-MW002-160407 4/7/2016	A11-MW002 E3Y50 A11-MW002-170310 3/10/2017	A11-MW002 E3YA3 A11-MW002-181114 11/14/2018	A11-MW002 E3YF7 A11-MW002-190521 5/21/2019	A11-MW002  A11-MW002-191113 11/13/2019	A11-MW002 A11-MW002-200304 A11-MW002-200304 3/4/2020	A11-MW002 E3YG8 A11-MW002-200610 6/10/2020	A11-MW002 A11-MW002-200910 A11-MW002-200910 9/10/2020	A11-MW002 A11-MW002-201201 A11-MW002-201201 12/2/2020
Analyte Name	RG									
1,1,1-Trichloroethane	200	190 J	1000 U	250 U	2500 U	200 U	100 U	1 J	50.0 U	50.0 U
1,1,2-Trichloroethane	5	500 U	1000 U	250 U	2500 U	200 U	100 U	5 U	50.0 U	50.0 U
1,1-Dichloroethane	1400	500 U	1000 U	250 U	2500 U	200 U	100 U	6.4	50.0 U	50.0 U
1,1-Dichloroethene	7	500 U	1000 UJ	250 U	2500 U	200 U	100 U	5 U	50.0 U	50.0 U
1,2,4-Trimethylbenzene	--	NA	NA	NA	NA	403	822	NA	622	588
1,2-Dichlorobenzene	600	500 U	1000 U	250 U	2500 U	200 U	100 U	5	50.0 U	50.0 U
1,2-Dichloropropane	5	500 U	1000 U	250 U	2500 U	200 U	100 U	5 U	50.0 U	50.0 U
1,3,5-Trimethylbenzene	--	NA	NA	NA	NA	200 U	285	NA	202	161
1,4-Dioxane	7.7	NA	NA	NA	NA	4.42	3.31	4.03	2.9	1.1
2-Butanone	4200	1000 U	2000 U	500 U	5000 U	NA	625 U	10 U	312 U	312 U
2-Hexanone	--	1000 U	2000 U	500 U	5000 U	NA	250 U	10 U	125 U	125 U
4-Methyl 2-Pentanone	--	1000 U	2000 U	500 U	5000 U	NA	250 U	10 U		125 U
Acetone	6300	1000 U	2000 U	500 U	5000 U	NA	625 UJ	2.7 J	312 U	312 U
Benzene	5	500 U	1000 U	250 U	2500 U	200 U	100 U	5 U	50.0 U	50.0 U
Chloroethane	--	500 U	1000 U	250 U	2500 U	200 U	100 U	5 U	50.0 U	50.0 U
cis-1,2-Dichloroethene	70	140 J	1000 UJ	170 J	2500 U	200 U	100 U	32	50.0 U	50.0 U
Cyclohexane	--	500 U	1000 U	110 J	2500 U	NA	NA	120	NA	NA
Dichlorodifluoromethane (Freon 12)	1400	500 U	1000 U	250 U	2500 U	200 UJ	100 U	5 U	50.0 UJ	50.0 U
Ethyl Benzene	700	5700	4100	3300	3700	4420 J	6840	6400	8260	10200
Isopropyl Benzene	700	500 U	1000 U	93 J	2500 U	200 U	121	98	90	78.1
Methyl Acetate	--	500 U	1000 U	250 U	2500 U	NA	NA	11	NA	NA
Methylcyclohexane	--	440 J	530 J	780	780 J	NA	NA	570 J	NA	NA
Naphthalene	140	NA	NA	NA	NA	200 U	100 U	NA	55.2	58.5
n-Propylbenzene	--	NA	NA	NA	NA	200 U	215	NA	129	87.4
Styrene	100	500 U	1000 U	250 U	2500 U	371	100 U	5 U	50.0 U	50.0 U
Tetrachloroethene	5	500 U	1000 U	250 U	2500 U	200 U	100 U	5 U	50.0 U	50.0 U
Toluene	1000	180000 J	220000 J	160000 J	88000	22500 J	78600	68000 J	39300	33200
Trichloroethene	5	500 U	1000 U	250 U	2500 U	200 U	100 U	0.86 J	50.0 U	50.0 U
Vinyl Chloride	2	500 U	1000 U	250 U	2500 U	200 U	100 U	4.4 J	50.0 U	50.0 U
Xylenes (Total)	10000	25200	17800	3300	14200	14930 J	24900	24900	32820	38040

Notes:  
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Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW003 E52K8 A11-MW003-110420 4/20/2011	A11-MW003 E52N6 A11-MW003-110721 7/21/2011	A11-MW003 E52S3 A11-MW003-1201112 1/12/2012	A11-MW003 E52S9 A11-MW003-04/03/2012 4/3/2012	A11-MW003 E3XC1 A11-MW003-120919 9/19/2012	A11-MW003 E3XD1 A11-MW003-121205 12/5/2012	A11-MW003 E3XG0 A11-MW003-130314 3/14/2013	A11-MW003 E3XG9 A11-MW003-130626 6/26/2013	A11-MW003 E3XP8 A11-MW003-140807 8/7/2014	A11-MW003 E3XQ8 A11-MW003-141217 12/17/2014	A11-MW003 E3XX6 A11-MW003-150520 5/20/2015
Analyte Name	RG											
1,1,1-Trichloroethane	200	1000 U	27	200 U	50 U	5000 U	130 U	130 U	7.4	130 U	50 U	50 UJ
1,1-Dichloroethane	1400	1000 U	20	200 U	11 J	5000 U	130 U	130 U	9.2	130 U	5.5 J	4.8 J-
1,1-Dichloroethene	7	130 J	10 U	200 U	50 U	5000 U	130 U	130 U	5 U	130 U	50 U	50 UJ
1,2,4-Trimethylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,3,5-Trimethylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,4-Dioxane	7.7	9300 J	200 R	4000 R	1000 R	100000 R	2500 U	2500 R	100 R	2500 R	1000 R	1000 R
2-Chlorotoluene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
4-Methyl 2-Pentanone	--	2000 U	20 U	400 U	100 U	10000 U	250 U	250 U	10 U	250 U	100 U	100 UJ
Acetone	6300	4000 U	40 U	400 U	100 U	10000 U	250 U	250 U	10 U	250 U	100 U	100 UJ
Chloroethane	--	1000 U	10 U	200 U	50 U	5000 U	130 U	130 U	4.2 J	130 U	50 U	50 UJ
cis-1,2-Dichloroethene	70	1000 U	2.7 J	200 U	50 U	5000 U	130 U	130 U	1.3 J	130 U	50 U	50 UJ
Cyclohexane	--	1000 U	4.7 J	200 U	10 J	5000 U	130 U	130 U	5 U	130 U	50 U	50 UJ
Ethyl Benzene	700	1200	420 D	3000	1300	4500 J	300	92 J	40 J	730	78	320 J-
Isopropyl Benzene	700	1000 U	31	85 J	66	5000 U	130 U	53 J	58	130 U	36 J	51 J-
Methylcyclohexane	--	1000 U	35	160 J	91	5000 U	140	140	160	380	250	50 UJ
Methylene Chloride	5	1000 U	0.65 J	200 U	100 U	5000 U	5 U	130 U	5 U	130 U	50 U	100 UJ
Naphthalene	140	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
n-Butylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
n-Propylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
sec-Butylbenzene	--	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Toluene	1000	1000	10 U	1000	550	130000	980	130 U	32 J	620	96	160 J-
Trichloroethene	5	1000 U	10 U	200 U	50 U	5000 U	130 U	130 U	0.53 J	130 U	50 U	50 UJ
Xylenes (Total)	10000	5400 J	3003	12200 D	11025 D	16800	9830	4200	1400 D	3200	1200	4300 DJ-

Notes:  
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Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW003 E3Y00 A11-MW003-160407 4/7/2016	A11-MW003 E3Y51 A11-MW003-170310 3/10/2017	A11-MW003 E3YA4 A11-MW003-181114 11/14/2018	A11-MW003 E3YF4 A11-MW003-190521 5/21/2019	A11-MW003  A11-MW003-191113 11/13/2019	A11-MW003 A11-MW003-200304 A11-MW003-200304 3/4/2020	A11-MW003 E3YG5 A11-MW003-200610 6/10/2020	A11-MW003 A11-MW003-200910 A11-MW003-200910 9/10/2020	A11-MW003 A11-MW003-201201 A11-MW003-201201 12/2/2020
Analyte Name	RG									
1,1,1-Trichloroethane	200	2.3 J	250 U	100 U	250 U	10.0 U	20.0 U	1.3 J	10.0 U	10.0 U
1,1-Dichloroethane	1400	5.3	250 U	100 U	250 U	10.0 U	20.0 U	6.9	10.0 U	10.0 U
1,1-Dichloroethene	7	5 UJ	250 U	100 U	250 U	10.0 U	20.0 U	5 U	10.0 U	10.0 U
1,2,4-Trimethylbenzene	--	NA	NA	NA	NA	137	329	NA	113	178
1,3,5-Trimethylbenzene	--	NA	NA	NA	NA	47.1	121	NA	34.8	55
1,4-Dioxane	7.7	NA	NA	NA	NA	12	8.57	9.58	7.23	4.6
2-Chlorotoluene	--	NA	NA	NA	NA	10.3	20.0 U	NA	10.0 U	10.0 U
4-Methyl 2-Pentanone	--	10 U	500 U	200 U	500 U	NA	50.0 U	2.2 J	25.0 U	25.0 U
Acetone	6300	10 U	500 U	200 U	78 J	NA	125 UJ	10 U	62.5 U	62.5 U
Chloroethane	--	5 U	250 U	100 U	250 U	10.0 U	20.0 U	1.9 J	10.0 U	10.0 U
cis-1,2-Dichloroethene	70	1.2 J-	250 U	100 U	250 U	10.0 U	20.0 U	5 U	10.0 U	10.0 U
Cyclohexane	--	1.8 J	250 U	100 U	250 U	NA	NA	7.8	NA	NA
Ethyl Benzene	700	33	450	130	160 J	144	1500	430 J	201	256
Isopropyl Benzene	700	6.4	250 U	100 U	57 J	31.7	74.4	45	28.3	38.5
Methylcyclohexane	--	43	250 U	100 U	110 J	NA	NA	59	NA	NA
Methylene Chloride	5	5 U	250 U	100 U	250 U	10.0 U	20.0 U	5 U	10.0 U	10.0 U
Naphthalene	140	NA	NA	NA	NA	13.8	29.1	NA	10.0 U	16.6
n-Butylbenzene	--	NA	NA	NA	NA	10.0 U	23.7	NA	10.0 U	13.1
n-Propylbenzene	--	NA	NA	NA	NA	33.6	76	NA	28.2	37.3
sec-Butylbenzene	--	NA	NA	NA	NA	10.0 U	27.8	NA	12	15
Toluene	1000	23	190 J	54 J	570	133	38.4	7.6 J	10.0 U	10.0 U
Trichloroethene	5	5 U	250 U	100 U	250 U	10.0 U	20.0 U	5 U	10.0 U	10.0 U
Xylenes (Total)	10000	392.2 J	4900	3500	12000	2910	13000	5105.2	2430	6310

Station Location EPA Sample ID Sample ID Sample Date		A11-MW004A E52K9 A11-MW004A-110420 4/20/2011	A11-MW004A E52N1 A11-MW004A-110720 7/20/2011	A11-MW004A E52S2 A11-MW004A-120112 1/12/2012	A11-MW004A E52T0 A11-MW004A-04/03/2012 4/3/2012	A11-MW004A E3XC2 A11-MW004A-120919 9/19/2012	A11-MW004A E3XD2 A11-MW004A-121205 12/5/2012	A11-MW004A E3XG1 A11-MW004A-130314 3/14/2013	A11-MW004A E3XH1 A11-MW004A-130626 6/26/2013	A11-MW004A E3XP9 A11-MW004A-140807 8/7/2014	A11-MW004A E3XQ9 A11-MW004A-141217 12/17/2014	A11-MW004A E3XX7 A11-MW004A-150520 5/20/2015
Analyte Name	RG											
1,1,1-Trichloroethane	200	10000 U	1000 U	2500 U	40 J	5000 U	1000 U	500 U	42	5000 U	1000 U	1000 UJ
1,1-Dichloroethane	1400	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	2 J	5000 U	1000 U	1000 UJ
1,1-Dichloroethene	7	1100 J	1000 U	2500 U	250 U	5000 U	1000 U	500 U	9.4	5000 U	1000 U	1000 UJ
1,4-Dioxane	7.7	200000 R	20000 R	50000 R	5000 R	100000 R	20000 U	10000 R	100 R	100000 R	20000 R	20000 R
2-Butanone	4200	20000 U	2000 U	5000 U	500 U	10000 U	2000 U	1000 U	1.2 J	10000 U	2000 U	2000 UJ
Acetone	6300	40000 U	4000 U	5000 U	500 U	10000 U	2000 U	1000 U	2 J	10000 U	2000 U	2000 UJ
cis-1,2-Dichloroethene	70	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	17	5000 U	1000 U	1000 UJ
Cyclohexane	--	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	5 U	5000 U	1000 U	1000 UJ
Ethyl Benzene	700	10000 U	240 J	3000	400	5000 U	430 J	1100	810 D	5000 U	220 J	420 J-
Isopropyl Benzene	700	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	7.1 J	5000 U	1000 U	1000 UJ
Methylcyclohexane	--	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	28 J	5000 U	1000 U	1000 UJ
Styrene	100	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	5 UJ	5000 U	1000 U	1000 UJ
Tetrachloroethene	5	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	37 J	5000 U	1000 U	1000 UJ
Toluene	1000	160000	200000 D	180000 D	120000 D	170000	120000 D	190000 D	230000 D	110000	100000 D	130000 J-D
Trichloroethene	5	10000 U	1000 U	2500 U	250 U	5000 U	1000 U	500 U	5.8 J	5000 U	1000 U	1000 UJ
Xylenes (Total)	10000	10000 U	419 J	12400	707 J	2100 J	2250 J	4570	3660	3900 J	430 J	98 J-

Notes:  
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Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW004A E3Y01 A11-MW004A-160407 4/7/2016	A11-MW004A E3Y52 A11-MW004A-170310 3/10/2017	A11-MW004A E3YA6 MW004A-181115 11/14/2018	A11-MW004A** E3YA7 MW004A-181115-D 11/14/2018	A11-MW004A E3YF8 A11-MW004A-190521 5/21/2019	A11-MW004A  A11-MW004A-191113 11/13/2019	A11-MW004A A11-MW004A-200304 A11-MW004A-200304 3/4/2020	A11-MW004A E3YG9 A11-MW004A-200610 6/10/2020	A11-MW004A A11-MW004A-200910 A11-MW004A-200910 9/10/2020	A11-MW004A A11-MW004A-201201 A11-MW004A-201201 12/2/2020
Analyte Name	RG										
1,1,1-Trichloroethane	200	21	250 U	2500 U	2500 U	250 U	400 U	50.0 U	3.7 J	50.0 U	50.0 U
1,1-Dichloroethane	1400	1.3 J	250 U	2500 U	2500 U	250 U	400 U	50.0 U	5 U	50.0 U	50.0 U
1,1-Dichloroethene	7	5 U	250 UJ	2500 UJ	2500 UJ	250 U	400 U	50.0 U	5 U	50.0 U	50.0 U
1,4-Dioxane	7.7	NA	NA	NA	NA	NA	1.2	1.41	1.51	1.09	NA
2-Butanone	4200	10 U	500 U	25000 U	25000 U	500 U	NA	312 U	10 U	312 U	312 U
Acetone	6300	10 U	500 U	25000 U	25000 U	99 J	NA	312 UJ	10 U	312 U	312 U
cis-1,2-Dichloroethene	70	14	250 UJ	2500 UJ	2500 UJ	250 U	400 U	50.0 U	5 U	50.0 U	50.0 U
Cyclohexane	--	5 U	250 U	2500 U	2500 U	250 U	NA	NA	0.83 J	NA	NA
Ethyl Benzene	700	440 J	320	2500 U	2500 U	440	455	260	330 J	365	331
Isopropyl Benzene	700	4 J-	250 U	2500 U	2500 U	250 U	400 U	50.0 U	2.1 J	50.0 U	50.0 U
Methylcyclohexane	--	22	250 U	2500 U	2500 U	250 U	NA	NA	21	NA	NA
Styrene	100	5 UJ	250 U	2500 U	2500 U	250 U	400 U	50.0 U	5 U	50.0 U	50.0 U
Tetrachloroethene	5	18 J-	250 U	2500 U	2500 U	250 U	400 U	50.0 U	5.3	50.0 U	50.0 U
Toluene	1000	150000 J	79000	48000	39000	59000	64300	45300	52000	42600	34200
Trichloroethene	5	8.8 J-	250 U	2500 U	2500 U	250 U	400 U	50.0 U	1 J	50.0 U	50.0 U
Xylenes (Total)	10000	1140 D	539 J	2500 U	2500 U	706 J	800 U	414	531 J	604.6	541.2

Station Location EPA Sample ID Sample ID Sample Date		A11-MW004B E52L0 A11-MW004B-110420 4/20/2011	A11-MW004B E52N2 A11-MW004B-110720 7/20/2011	A11-MW004B E52S1 A11-MW004B-120112 1/12/2012	A11-MW004B E52T1 A11-MW004B-04/03/2012 4/3/2012	A11-MW004B** E52T6 A11-MW004B-04/03/2012D 4/3/2012	A11-MW004B E3XC3 A11-MW004B-120919 9/18/2012	A11-MW004B** E3XC4 A11-MW004B-120919-D 9/18/2012	A11-MW004B E3XD3 A11-MW004B-121204 12/4/2012	A11-MW004B** E3XD4 A11-MW004B-121204-D 12/4/2012	A11-MW004B E3XG2 A11-MW004B-130314 3/14/2013	A11-MW004B** E3XG3 A11-MW004B-130314-D 3/14/2013
Analyte Name	RG											
1,1,1-Trichloroethane	200	190 J	98	74	59	58	64	64	26	38	24 D	35 D
1,1,2-Trichloroethane	5	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 U	5 U	0.5 U	0.5 U
1,1-Dichloroethane	1400	20 J	13	11	9	9.1	8.2	8.5	6.7	8.3	5.9	5.9
1,1-Dichloroethene	7	6.6 J	5.3	5 U	3.7 J	4.0 J	5 U	5 U	5 U	5 U	2.1	2.2
1,4-Dioxane	7.7	400 R	100 R	100 R	100 R	100 R	100 R	100 R	100 U	100 U	NA	NA
Acetone	6300	80 UJ	20 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	5 U	5 U
Benzene	5	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 UJ	5 U	0.5 U	0.5 U
cis-1,2-Dichloroethene	70	7.8 J	5.1	5.9	4.1 J	4.0 J	4.2 J	4.4 J	3.3 J	3.9 J	2.9	2.9
Dichlorodifluoromethane (Freon 12)	1400	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 U	5 U	0.5 U	0.5 U
Ethyl Benzene	700	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 U	5 U	0.15 J	0.16 J
Tetrachloroethene	5	20 UJ	0.4 J	5 U	0.67 J	0.55 J	0.61 J	0.79 J	5 U	5 U	0.39 J	0.36 J
Toluene	1000	20 UJ	5 U	5 U	3.0 J	2.8 J	0.75 J	0.94 J	5 U	2.1 J	0.5 U	0.5 U
trans-1,2-Dichloroethene	100	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 U	5 U	0.13 J	0.13 J
trans-1,3-Dichloropropene	--	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 U	5 U	0.5 U	0.5 U
Trichloroethene	5	4 J	2.7 J	3.4 J	3.0 J	2.9 J	3.5 J	3.6 J	1.9 J	2.4 J	1.9	1.8
Trichlorofluoromethane (Freon 11)	2100	20 UJ	5 U	5 U	5.0 U	5.0 U	5 U	5 U	5 U	5 U	0.21 J	0.18 J
Xylenes (Total)	10000	20 UJ	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	0.69 J	0.67 J

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Table 6  
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Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW004B E3XH2 A11-MW004B-130626 6/26/2013	A11-MW004B** E3XH3 A11-MW004B-130626-D 6/26/2013	A11-MW004B E3XP3 A11-MW004B-140807 8/7/2014	A11-MW004B** E3XP4 A11-MW004B-140807-D 8/7/2014	A11-MW004B E3XQ4RE A11-MW004B-141217RE 12/17/2014	A11-MW004B** E3XQ5RE A11-MW004B-141217-DRE 12/17/2014	A11-MW004B E3XX2 A11-MW004B-150520 5/20/2015	A11-MW004B** E3XX3 A11-MW004B-150520-D 5/20/2015	A11-MW004B E3Y02 A11-MW004B-160406 4/6/2016	A11-MW004B** E3Y03 A11-MW004B-160406-D 4/6/2016	A11-MW004B E3Y48 A11-MW004B-170310 3/10/2017
Analyte Name	RG											
1,1,1-Trichloroethane	200	25 D	27 D	14	14	15 J	16	14 J-	15 J-	12	11	12
1,1,2-Trichloroethane	5	0.066 J	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.5 U
1,1-Dichloroethane	1400	7.1	7.6	6.3	6.2	8.7	9.4	9 J-	9.2 J-	8.8	9.8	10
1,1-Dichloroethene	7	2.5	2.8	1.2 J	1.2 J	2.5	2	1.6 J-	1.6 J-	0.5 U	0.5 U	1.8
1,4-Dioxane	7.7	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Acetone	6300	10 U	10 U	5 U	5 U	10 U	10 U	5 UJ	5 UJ	5 U	5 U	5 U
Benzene	5	0.086 J	0.078 J	0.5 U	0.5 U	0.5 U	0.5 U	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.5 U
cis-1,2-Dichloroethene	70	3	3.2	1.8 J	1.7 J	2.2	2.2	2.1 J-	2.1 J-	2	2.2	2.1
Dichlorodifluoromethane (Freon 12)	1400	2.2	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.5 U
Ethyl Benzene	700	0.16 J	0.17 J	0.2 J	0.19 J	0.5 U	0.5 U	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.5 U
Tetrachloroethene	5	0.64	0.67	0.49 J	0.45 J	0.61	0.59	0.53 J-	0.53 J-	0.57	0.47 J	0.55
Toluene	1000	0.8	0.84	1 U	1 U	590 U	590 U	1.4 UJ	1.4 UJ	8	7.9	0.1 J
trans-1,2-Dichloroethene	100	0.21 J	0.21 J	0.16 J	0.13 J	0.24 J	0.25 J	0.26 J-	0.25 J-	0.25 J	0.22 J	0.23 J
trans-1,3-Dichloropropene	--	0.18 J	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.5 U
Trichloroethene	5	2.6	2.9	2.1	2.2	2	2	2.1 J-	2.2 J-	2.6	2.3	2
Trichlorofluoromethane (Freon 11)	2100	0.36 J	0.37 J	0.33 J	0.33 J	0.52 J	0.43 J	0.53 J-	0.51 J-	0.38 J	0.36 J	0.5 U
Xylenes (Total)	10000	1.01 J	1.1 J	0.31 J	0.3 J	110 U	110 U	0.5 UJ	0.5 UJ	0.38 J	0.14 J	0.11 J

Station Location EPA Sample ID Sample ID Sample Date		A11-MW004B E3YA5 A11-MW004B-181114 11/15/2018	A11-MW004B E3YF2 A11-MW004B-190520 5/20/2019	A11-MW004B A11-MW004B-191112 11/12/2019	A11-MW004B A11-MW004B-200303 A11-MW004B-200303 3/3/2020	A11-MW004B E3YG3 A11-MW004B-200609 6/9/2020	A11-MW004B A11-MW004B-200909 A11-MW004B-200909 9/9/2020	A11-MW004B E3YJ0 A11-MW004B-201201 12/1/2020
Analyte Name	RG							
1,1,1-Trichloroethane	200	8.9	18	6.26	5.29 J	5.5	4.93 J	5.61
1,1,2-Trichloroethane	5	5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
1,1-Dichloroethane	1400	11	20	6.55	5.86	6.3	5.34	5.67
1,1-Dichloroethene	7	5 U	3.2	2.00 U	2.00 U	0.95	2.00 U	2.00 U
1,4-Dioxane	7.7	NA	NA	13.4	9.75	11.7	7.86	6.3
Acetone	6300	4.5 J	6.1 U	NA	12.5 UJ	5 U	12.5 U	12.5 U
Benzene	5	5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
cis-1,2-Dichloroethene	70	1.8 J	3.8	2.00 U	2.00 U	1.3	2.00 U	2.00 U
Dichlorodifluoromethane (Freon 12)	1400	5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Ethyl Benzene	700	5 U	0.5 UJ	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Tetrachloroethene	5	5 U	0.93 J-	2.00 U	2.00 U	0.4 J	2.00 U	2.00 U
Toluene	1000	1.6 J	0.5 UJ	2.00 U	2.00 U	1.6	2.00 UJ	2.00 U
trans-1,2-Dichloroethene	100	5 U	0.52	2.00 U	2.00 U	0.16 J	2.00 U	2.00 U
trans-1,3-Dichloropropene	--	5 U	0.5 U	2.00 U	2.00 UJ	0.5 U	2.00 U	2.00 U
Trichloroethene	5	1.6 J	3.1 J-	2.00 U	2.00 U	1.4	2.00 U	2.00 U
Trichlorofluoromethane (Freon 11)	2100	5 U	0.29 J	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Xylenes (Total)	10000	5 U	0.77 J	4.00 U	4.00 U	0.5 U	4.00 U	4.00 U

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
U = Not detected at value shown  
J = Estimated result      J- = Estimated result biased low      J+ = Estimated result biased high  
R = Rejected



Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW005 E52H3 A11-MW005-110419 4/19/2011	A11-MW005** E52H4 A11-MW005-110419-D 4/19/2011	A11-MW005 E52L6 A11-MW005-110720 7/20/2011	A11-MW005** E52L7 A11-MW005-110720-D 7/20/2011	A11-MW005 E52S0 A11-MW005-120111 1/11/2012	A11-MW005 E52T2 A11-MW005-04/02/2012 4/2/2012	A11-MW005 E3XC5 A11-MW005-120918 9/18/2012	A11-MW005 E3XD5 A11-MW005-121204 12/4/2012	A11-MW005 E3XG4 A11-MW005-130313 3/13/2013	A11-MW005 E3XH4 A11-MW005-130625 6/25/2013	A11-MW005 E3XP1 A11-MW005-140806 8/6/2014
Analyte Name	RG											
1,1,1-Trichloroethane	200	67	69	38	38	15	16	19	12	13	11	3.6
1,1-Dichloroethane	1400	15	15	13	13	7	8.1	7.4	9.5	7.8	7.3	2.8
1,1-Dichloroethene	7	5 U	25 U	3.6 J	3.8 J	5 U	2.4 J	5 U	5 U	1.8	0.5 U	1 U
1,4-Dioxane	7.7	21 J	17 J	100 R	100 R	100 R	100 R	100 R	100 U	NA	NA	NA
Benzene	5	5 U	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.096 J	0.5 U
Bromodichloromethane	0.2*	5 U	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U
Chloroform	70	5 U	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U
cis-1,2-Dichloroethene	70	6	6	4.6 J	4.7 J	3.2 J	2.9 J	2.6 J	3.5 J	3.2	2.7	1.4 J
Dibromochloromethane	140*	5 U	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U
Dichlorodifluoromethane (Freon 12)	1400	2.5 J	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U
Ethyl Benzene	700	5 U	5 U	0.18 J	0.23 J	5 U	0.80 J	5 U	5 U	0.14 J	0.18 J	0.25 J
Isopropyl Benzene	700	5 U	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U
Tetrachloroethene	5	5 U	5 U	0.4 J	0.41 J	5 U	0.45 J	5 U	5 U	0.29 J	0.41 J	0.23 J
Toluene	1000	5 U	5 U	5 U	5 U	5 U	1.8 J	0.66 J	5 U	0.5 U	0.82	1 U
trans-1,2-Dichloroethene	100	5 U	5 U	0.5 J	5 U	5 U	5.0 U	5 U	5 U	0.18 J	0.19 J	0.5 U
Trichloroethene	5	1.8 J	1.7 J	1.4 J	1.4 J	5 U	0.95 J	1.2 J	5 U	0.97	1.3 J	0.5
Trichlorofluoromethane (Freon 11)	2100	5 U	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U
Xylenes (Total)	10000	5 U	5 U	0.31 J	0.49 J	5 U	0.42 J	5 U	5 U	0.43 J	0.92 J	0.35 J

Station Location EPA Sample ID Sample ID Sample Date		A11-MW005 E3XQ2 A11-MW005-141217 12/17/2014	A11-MW005 E3XX0 A11-MW005-150519 5/19/2015	A11-MW005 E3Y04 A11-MW005-160406 4/6/2016	A11-MW005 E3YA8 A11-MW005-181113 11/13/2018	A11-MW005 E3YF0 A11-MW005-190520 5/20/2019	A11-MW005 A11-MW005-191112 11/12/2019	A11-MW005 A11-MW005-200303 A11-MW005-200303 3/3/2020	A11-MW005 E3YG1 A11-MW005-200609 6/9/2020	A11-MW005 A11-MW005-200909 A11-MW005-200909 9/9/2020	A11-MW005 E3YH9 A11-MW005-201201 12/1/2020
Analyte Name	RG										
1,1,1-Trichloroethane	200	5.4	6.1 J-	4.4	3.5	6.4	2.46	2.92 J	4.5	5.56 J	4.9
1,1-Dichloroethane	1400	3.6	2.9 J-	2.8	3.8	7.2	3.21	3.77	6.4	9.11	7.01
1,1-Dichloroethene	7	1.3	0.88 J-	0.5 U	0.5 UJ	1.1	2.00 U	2.00 U	1.1	2.00 UJ	2.00 U
1,4-Dioxane	7.7	NA	NA	NA	NA	NA	7.63	5.35	8.83	8.18 J	4.5 J
Benzene	5	0.5 U	0.5 UJ	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Bromodichloromethane	0.2*	0.5 UJ	0.5 UJ	0.12 J	0.12 J	0.33 J	2.00 U	2.00 U	0.4 J	2.00 U	2.00 U
Chloroform	70	0.5 U	0.5 UJ	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.45 J	2.00 U	2.00 U
cis-1,2-Dichloroethene	70	2.1	2.1 J-	1.9	0.85 J-	1.6	2.00 U	2.00 U	1.3	2.00 U	2.00 U
Dibromochloromethane	140*	0.5 U	0.5 UJ	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.18 J	2.00 U	2.00 U
Dichlorodifluoromethane (Freon 12)	1400	0.5 U	0.5 UJ	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Ethyl Benzene	700	0.5 UJ	0.5 UJ	0.21 J	0.5 UJ	0.14 J	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Isopropyl Benzene	700	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.14 J+	2.00 U	2.00 U	0.5 U	2.00 UJ	2.00 U
Tetrachloroethene	5	0.33 J	0.5 UJ	0.22 J	0.38 J	0.86	2.00 U	2.00 U	0.39 J	2.00 UJ	2.00 U
Toluene	1000	0.5 UJ	0.5 UJ	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
trans-1,2-Dichloroethene	100	0.5 U	0.11 J-	0.5 U	0.5 UJ	0.3 J	2.00 U	2.00 U	0.15 J	2.00 UJ	2.00 U
Trichloroethene	5	1.2 J	1.7 J-	1.2	1.4	2.1	2.00 U	2.00 U	0.89	2.00 UJ	2.00 U
Trichlorofluoromethane (Freon 11)	2100	0.5 U	0.5 UJ	0.5 U	0.5 U	0.31 J	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Xylenes (Total)	10000	110 UJ	0.5 UJ	1.27 J	0.5 U	1.19 J	4.00 U	4.00 U	0.5 U	4.00 U	4.00 U

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
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Table 6  
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Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW006 E52L1 A11-MW006-110420 4/20/2011	A11-MW006 E52N3 A11-MW006-110720 7/20/2011	A11-MW006 E52R8 A11-MW006-120111 1/11/2012	A11-MW006** E52R9 A11-MW006-120111-D 1/11/2012	A11-MW006 E52T3 A11-MW006-04/02/2012 4/2/2012	A11-MW006 E3XC6 A11-MW006-120918 9/18/2012	A11-MW006 E3XD6 A11-MW006-121204 12/4/2012	A11-MW006 E3XG5 A11-MW006-130313 3/13/2013	A11-MW006 E3XH5RE A11-MW006-130625RE 6/25/2013	A11-MW006 E3XP0 A11-MW006-140806 8/6/2014	A11-MW006 E3XQ1 A11-MW006-141217 12/17/2014
Analyte Name	RG											
1,1,1-Trichloroethane	200	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 UJ
1,1-Dichloroethane	1400	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.17 J	0.11 J	0.5 U	0.12 J
1,1-Dichloroethene	7	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 UJ	0.5 U	0.5 U	0.5 UJ
1,2-Dichloropropane	5	5 U	0.73 J	5 U	5 U	5.0 U	5 U	5 U	1.2	0.62	0.5 U	0.5 UJ
1,4-Dioxane	7.7	100 R	100 R	100 R	100 R	100 R	100 R	100 U	NA	NA	NA	NA
Benzene	5	3.3 J	2.9 J	5 U	5 U	3.1 J	5 U	5 U	4.3	0.4 J	0.58	1.2 J
Bromochloromethane	--	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 UJ
Bromodichloromethane	0.2*	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 UJ
Chloroethane	--	3.8 J	5 U	5 U	5 U	1.2 J	0.93 J	5 U	4.7	0.3 J	0.5 U	0.5 UJ
Chloroform	70	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 UJ
cis-1,2-Dichloroethene	70	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 UJ	0.5 U	0.5 U	0.1 J
Cyclohexane	--	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	1	1.2 J
Dibromochloromethane	140*	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.5 U	0.5 UJ
Dichlorodifluoromethane (Freon 12)	1400	2 J	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.088 J	0.5 U	0.5 UJ
Ethyl Benzene	700	5 U	0.21 J	5 U	5 U	5.0 U	5 U	5 U	0.21 J	0.21 J	0.36 J	0.5 UJ
Isopropyl Benzene	700	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 UJ	0.5 U	6.3	3.7 J
Methylcyclohexane	--	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 U	0.5 U	0.32 J	0.5 UJ
Tetrachloroethene	5	1.2 J	5 U	5 U	5 U	0.63 J	1.4 J	5 U	0.46 J	1	0.38 J	0.53 J
Toluene	1000	5 U	5 U	5 U	5 U	5.0 U	0.89 J	5 U	0.5 UJ	0.9	1 U	0.5 UJ
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 UJ	0.5 U	0.5 U	0.5 UJ
Trichloroethene	5	5 U	5 U	5 U	5 U	5.0 U	5 U	5 U	0.5 UJ	0.5 U	0.5 U	0.5 UJ
Xylenes (Total)	10000	1.3 J	0.71 J	5 U	5 U	5.0 U	5 U	5 U	0.8 J	1.23 J	1.66 J	110 UJ

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
U = Not detected at value shown  
J = Estimated result      J- = Estimated result biased low      J+ = Estimated result biased high  
R = Rejected



Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW006 E3XW9 A11-MW006-150519 5/19/2015	A11-MW006 E3Y05 A11-MW006-160406 4/6/2016	A11-MW006 E3Y47 A11-MW006-170309 3/9/2017	A11-MW006 E3YA9 A11-MW006-181113 11/13/2018	A11-MW006 E3YE9 A11-MW006-190520 5/20/2019	A11-MW006  A11-MW006-191113 11/13/2019	A11-MW006 A11-MW006-200303 A11-MW006-200303 3/3/2020	A11-MW006 E3YG0 A11-MW006-200609 6/9/2020	A11-MW006 A11-MW006-200909 A11-MW006-200909 9/9/2020	A11-MW006 E3YH8 A11-MW006-201201 12/1/2020
Analyte Name	RG										
1,1,1-Trichloroethane	200	0.5 UJ	0.23 J	0.38 J	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
1,1-Dichloroethane	1400	0.5 UJ	0.72	1.8	0.2 J	0.75	2.00 U	2.00 U	0.11 J	2.00 U	2.00 U
1,1-Dichloroethene	7	0.5 UJ	0.5 UJ	0.5 U	0.5 UJ	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
1,2-Dichloropropane	5	0.5 UJ	0.5 U	0.5 U	0.5 U	3.1	2.57	2.00 U	0.5 U	2.00 U	2.00 U
1,4-Dioxane	7.7	NA	NA	NA	NA	NA	1.02	1.54	7.53	8.42	4.1
Benzene	5	0.5 UJ	0.5 U	0.5 U	2.5	8.8	3.12	2.62	2	2.28	2.82
Bromochloromethane	--	0.5 UJ	0.21 J	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Bromodichloromethane	0.2*	0.5 UJ	0.39 J	0.75	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Chloroethane	--	0.5 UJ	0.5 U	0.5 U	2.3	1.4	2.00 U	2.00 U	0.44 J	2.00 U	2.00 U
Chloroform	70	0.5 UJ	0.5 U	0.6	0.5 U	0.5 U	2.00 U	2.00 U	2.00 U	2.00 U	2.00 U
cis-1,2-Dichloroethene	70	0.5 UJ	0.47 J-	0.31 J	0.5 UJ	0.74 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Cyclohexane	--	0.5 UJ	0.5 U	0.5 U	0.5 U	0.5 U	NA	NA	0.35 J	NA	NA
Dibromochloromethane	140*	0.5 UJ	0.45 J	0.51	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Dichlorodifluoromethane (Freon 12)	1400	0.5 UJ	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Ethyl Benzene	700	0.5 UJ	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Isopropyl Benzene	700	1.6 J-	0.5 U	0.5 U	22	37 J+	2.00 U	2.00 U	0.14 J	2.00 U	2.00 U
Methylcyclohexane	--	0.5 UJ	0.5 U	0.5 U	0.5 U	0.71	NA	NA	0.5 U	NA	NA
Tetrachloroethene	5	0.54 J-	0.41 J	0.82	0.15 J	0.17 J	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Toluene	1000	1.4 UJ	0.5 U	0.5 U	0.5 U	0.5 U	2.00 U	2.00 U	0.5 U	2.00 UJ	2.00 U
trans-1,2-Dichloroethene	100	0.5 UJ	0.5 UJ	0.5 U	0.5 UJ	0.5 U	2.00 U	2.00 U	0.5 U	2.00 U	2.00 U
Trichloroethene	5	0.5 UJ	0.41 J	0.28 J	0.5 U	0.5 U	2.00 U	2.00 U	0.14 J	2.00 U	2.00 U
Xylenes (Total)	10000	0.5 UJ	0.13 J	0.5 U	0.5 U	0.24 J+	4.00 U	4.00 U	0.5 U	4.00 U	4.00 U

Station Location EPA Sample ID Sample ID Sample Date		A11-MW007 E3YB0 A11-MW007-181114 11/14/2018	A11-MW007** E3YB1 A11-MW007-181114-D 11/14/2018	A11-MW007 E3YF5 A11-MW007-190521 5/21/2019	A11-MW007** E3YF6 A11-MW007-190521-D 5/21/2019	A11-MW007  A11-MW007-191113 11/13/2019	A11-MW007**  A11-MW007-191113-D 11/13/2019	A11-MW007 A11-MW007-200304 A11-MW007-200304 3/4/2020	A11-MW007** A11-MW007-200304-D A11-MW007-200304-D 3/4/2020	A11-MW007 E3YG6 A11-MW007-200610 6/10/2020	A11-MW007** E3YG7 A11-MW007-200610-D 6/10/2020	A11-MW007 A11-MW007-200910 A11-MW007-200910 9/10/2020
Analyte Name	RG											
1,1-Dichloroethane	1400	250 U	250 U	20 J	21 J	10.0 U	10.0 U	4.00 U	4.00 U	5 U	5 U	10.0 U
1,2,4-Trimethylbenzene	--	NA	NA	NA	NA	31.7	32.1	22.6	21	NA	NA	53.5
1,3,5-Trimethylbenzene	--	NA	NA	NA	NA	10.0 U	10.0 U	4.84	4.57	NA	NA	11.1
1,4-Dioxane	7.7	NA	NA	NA	NA	0.278	0.293	3.38	3.3	0.205 U	0.205 U	0.212 U
Benzene	5	250 U	250 U	130 U	130 U	10.0 U	10.0 U	4.00 U	4.00 U	5 U	5 U	10.0 U
Ethyl Benzene	700	6500	6700	2500	2600	1420	1420	959	863	820	810	2630
Isopropyl Benzene	700	99 J	110 J	91 J	92 J	28.3	28.8	12	11	6.5	6.5	86.1
Methylcyclohexane	--	250 U	89 J	100 J	99 J	NA	NA	NA	NA	2.7 J	2.6 J	NA
Naphthalene	140	NA	NA	NA	NA	10.0 U	10.0 U	4.00 U	4.00 U	NA	NA	11.4
n-Butylbenzene	--	NA	NA	NA	NA	10.0 U	10.0 U	4.00 U	4.00 U	NA	NA	11.3
n-Propylbenzene	--	NA	NA	NA	NA	19.5	19.6	6.68	6.03	NA	NA	82.4
sec-Butylbenzene	--	NA	NA	NA	NA	10.0 U	10.0 U	4.00 U	4.00 U	NA	NA	10.8
Tetrachloroethene	5	250 U	250 U	130 U	130 U	10.0 U	10.0 U	4.00 U	4.00 U	1 J	0.89 J	10.0 U
Toluene	1000	200 J	230 J	12 J	10 J	10.0 U	10.0 U	4.00 U	4.00 U	5 UJ	5 UJ	10.0 U
Xylenes (Total)	10000	13000	13000	4500	4500	4190	4210	3050	2800	2600	2600	7600

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
U = Not detected at value shown  
J = Estimated result      J- = Estimated result biased low      J+ = Estimated result biased high  
R = Rejected



Table 6  
Comprehensive VOC Compounds Detected 2011-2020  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location EPA Sample ID Sample ID Sample Date		A11-MW007** A11-MW007-200910-D A11-MW007-200910-D 9/10/2020	A11-MW007 A11-MW007-201201 A11-MW007-201201 12/2/2020	A11-MW007** A11-MW007-201201-D A11-MW007-201201-D 12/2/2020
Analyte Name	RG			
1,1-Dichloroethane	1400	10.0 U	10.0 U	10.0 U
1,2,4-Trimethylbenzene	--	55.7	131	169
1,3,5-Trimethylbenzene	--	11.6	14.4 J	56.7 J
1,4-Dioxane	7.7	0.203 U	0.069 J	0.19 U
Benzene	5	10.0 U	10.0 UJ	44.3 J
Ethyl Benzene	700	2680	3300	3660
Isopropyl Benzene	700	89.1	109 J	486 J
Methylcyclohexane	--	NA	NA	NA
Naphthalene	140	13.2	34.0 J	97.3 J
n-Butylbenzene	--	12.4	19.9 J	66.9 J
n-Propylbenzene	--	84.7	104 J	454 J
sec-Butylbenzene	--	11.3	17.5 J	68.4 J
Tetrachloroethene	5	10.0 U	10.0 U	10.0 U
Toluene	1000	10.0 U	10.0 U	10.0 U
Xylenes (Total)	10000	7920	7390	8100

Station Location EPA Sample ID Sample ID Sample Date		A11-MW130A A11-MW130A-200909 A11-MW130A-200909 9/9/2020	A11-MW130A E3YH7 A11-MW130A-201201 12/1/2020
Analyte Name	RG		
1,1,1-Trichloroethane	200	3.51 J	3.51
1,1-Dichloroethane	1400	4.11	3.77
1,4-Dioxane	7.7	6.1	4
Total Xylenes	10000	4.00 U	4.00 U

Notes:  
All results in micrograms per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded results exceed remediation goal  
\*\* = Duplicate sample  
D = Diluted sample result  
U = Not detected at value shown  
J = Estimated result      J- = Estimated result biased low      J+ = Estimated result biased high  
R = Rejected



Table 7  
Comprehensive Attenuation Parameters  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location Sample ID Sample Date		A11-MW001 A11-MW001-140806 8/6/2014	A11-MW001 A11-MW-001-170309 3/9/2017	A11-MW001** A11-MW-001-170309-D 3/9/2017	A11-MW001 A11-MW001-181113 11/13/2018	A11-MW001 A11-MW001-190520 5/20/2019	A11-MW001 A11-MW001-191112 11/12/2019	A11-MW001 A11-MW001-200303 3/3/2020	A11-MW001 A11-MW001-200609 6/9/2020	A11-MW001 A11-MW001-200909 9/9/2020	A11-MW001 A11-MW001-201201 12/1/2020
Analyte Name	RG										
Alkalinity, Total (As CaCO3)	--	380	370	360	350	350	360	360	360	340	350
Methane	--	0.002 U	0.005 U	0.005 U	0.005 U	0.005 U	0.0027 J	0.00058 U	0.00058 U	0.00082 J	0.011
Nitrate	10	3.4 J	3 J-	3 J-	2.5	2.5	2.3	3.5	1.8 J	11.3	11.4
Sulfate	400	36.2	38.4	38.4	41.7	31	25	45	25	28.9 J-	30.4 J-

Station Location Sample ID Sample Date		A11-MW002 A11-MW002-140807 8/7/2014	A11-MW002 A11-MW-002-170310 3/10/2017	A11-MW002 A11-MW002-181114 11/14/2018	A11-MW002 A11-MW002-190521 5/21/2019	A11-MW002 A11-MW002-191113 11/13/2019	A11-MW002 A11-MW002-200304 3/4/2020	A11-MW002 A11-MW002-200610 6/10/2020	A11-MW002 A11-MW002-200910 9/10/2020	A11-MW002 A11-MW002-201201 12/2/2020
Analyte Name	RG									
Alkalinity, Total (As CaCO3)	--	350	390	380	370	400	370	400	420	440
Methane	--	2.9 (D)	15	13	20	26	16	19	26 J	31
Nitrate	10	0.2 U	0.2 R	0.2 U	1 U	0.2 U	0.2 U	0.2 UJ	0.12 U	0.12 U
Sulfate	400	10 U	35.6	10 U	20 U	4 U	4 U	4 U	0.12 U	1.09

Station Location Sample ID Sample Date		A11-MW003 A11-MW003-140807 8/7/2014	A11-MW003 A11-MW-003-170310 3/10/2017	A11-MW003 A11-MW003-181114 11/14/2018	A11-MW003 A11-MW003-190521 5/21/2019	A11-MW003 A11-MW003-191113 11/13/2019	A11-MW003** A11-MW003-191113-D 11/13/2019	A11-MW003 A11-MW003-200304 3/4/2020	A11-MW003 A11-MW003-200610 6/10/2020	A11-MW003 A11-MW003-200910 9/10/2020	A11-MW003 A11-MW003-201201 12/2/2020
Analyte Name	RG										
Alkalinity, Total (As CaCO3)	--	400	380	370	420	390	380	410	380	370	390
Methane	--	0.73 (D)	12	6.1	15	6.1	5.8	12	6.7	3.5 J	6.6
Nitrate	10	0.2 U	0.2 R	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 UJ	0.12 U	0.12 U
Sulfate	400	5.79	23.9	14.6	4 U	8.5	8.6	4 U	9	11.3	8.52

Station Location Sample ID Sample Date		A11-MW004A A11-MW004A-140807 8/7/2014	A11-MW004A A11-MW-004A-170310 3/10/2017	A11-MW004A A11-MW004A-181115 11/15/2018	A11-MW004A A11-MW004A-190521 5/21/2019	A11-MW004A A11-MW004A-191113 11/13/2019	A11-MW004A A11-MW004A-200304 3/4/2020	A11-MW004A** A11-MW004A-200304-D 3/4/2020	A11-MW004A A11-MW004A-200610 6/10/2020	A11-MW004A A11-MW004A-200910 9/10/2020	A11-MW004A A11-MW004A-201201 12/2/2020
Analyte Name	RG										
Alkalinity, Total (As CaCO3)	--	370	440	350	330	330	330	320	350	340	330
Methane	--	0.08 (D)	0.21	0.005 U	0.17	0.038	0.51	0.47	0.19	0.16 J	0.25
Nitrate	10	0.99	1.7 J-	1.5	1 U	0.23	0.28	0.38	0.65 J	1.93	1.66
Sulfate	400	9.48	93	26.9	30	32	36	36	35	33.4	42.9

Station Location Sample ID Sample Date		A11-MW004B A11-MW004B-140807 8/7/2014	A11-MW004B** A11-MW004B-140807-D 8/7/2014	A11-MW004B A11-MW-004B-170310 3/10/2017	A11-MW004B** A11-MW004B-181114-D 11/14/2018	A11-MW004B A11-MW004B-181114 11/15/2018	A11-MW004B A11-MW004B-190520 5/20/2019	A11-MW004B A11-MW004B-191112 11/12/2019	A11-MW004B A11-MW004B-200303 3/3/2020	A11-MW004B A11-MW004B-200609 6/9/2020	A11-MW004B A11-MW004B-200909 9/9/2020	A11-MW004B A11-MW004B-201201 12/1/2020
Analyte Name	RG											
Alkalinity, Total (As CaCO3)	--	370	350	340	360	350	350	350	330	340	340	340
Methane	--	0.002 U	0.002 U	0.005 U	0.04	0.046	0.005 U	0.00029 UJ	0.00058 U	0.035	0.027 J	0.02
Nitrate	10	1.8	1.8	1.8 J-	2.3	2.3	1.5	1.3	1.1	1.1 J	5.2	5.15
Sulfate	400	27.7	27.7	23.9	76.9	74.1	23	18	19	19	18.8 J-	18.9 J-

Notes:  
All results in milligram per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded result exceeds remediation goal  
\*\* = Duplicate sample  
U = Not detected above the reported limit  
J = Estimated result J- = Estimated result biased low J+ = Estimated result biased high  
(D) = Diluted result





Table 7  
Comprehensive Attenuation Parameters  
Source Area 11 2020 Groundwater Report  
Southeast Rockford Groundwater Contamination Superfund Site

Station Location Sample ID Sample Date		A11-MW005 A11-MW005-140806 8/6/2014	A11-MW005 A11-MW005-181113 11/13/2018	A11-MW005 A11-MW005-190520 5/20/2019	A11-MW005 A11-MW005-191112 11/12/2019	A11-MW005 A11-MW005-200303 3/3/2020	A11-MW005 A11-MW005-200609 6/9/2020	A11-MW005 A11-MW005-200909 9/9/2020	A11-MW005 A11-MW005-201201 12/1/2020
Analyte Name	RG								
Alkalinity, Total (As CaCO3)	--	380	360	370	370	320	370	370	380
Methane	--	0.002 U	0.005 U	0.005 U	0.00029 UJ	0.00058 U	0.00058 U	0.00094 J	0.00058 U
Nitrate	10	3.9 J	4.4	4.5	3.2	2.5	2.7 J	9.53	14.1
Sulfate	400	38.4	50.9	50	33	31	33	25.2 J-	33.7 J-

Station Location Sample ID Sample Date		A11-MW006 A11-MW006-140806 8/6/2014	A11-MW006 A11-MW-006-170309 3/9/2017	A11-MW006 A11-MW006-181113 11/13/2018	A11-MW006 A11-MW006-190520 5/20/2019	A11-MW006 A11-MW006-191112 11/12/2019	A11-MW006 A11-MW006-200303 3/3/2020	A11-MW006 A11-MW006-200609 6/9/2020	A11-MW006 A11-MW006-200909 9/9/2020	A11-MW006 A11-MW006-201201 12/1/2020
Analyte Name	RG									
Alkalinity, Total (As CaCO3)	--	10 U	340	400	450	490	490	460	440	450
Methane	--	0.64 (D)	0.18	5	6.5	3.4 J	3.3	3.8	4.1 J	8.1
Nitrate	10	0.2 UJ	2.1 J-	0.2 U	1 U	0.2 U	0.2 U	0.2 UJ	0.12 U	0.12 U
Sulfate	400	5 U	38.4	10 U	20 U	16	35	7.1	5.02 J-	5.56 J-

Station Location Sample ID Sample Date		A11-MW007 A11-MW007-181114 11/14/2018	A11-MW007 A11-MW007-190521 5/21/2019	A11-MW007** A11-MW007-190521-D 5/21/2019	A11-MW007 A11-MW007-191113 11/13/2019	A11-MW007 A11-MW007-200304 3/4/2020	A11-MW007 A11-MW007-200610 6/10/2020	A11-MW007** A11-MW007-200610-D 6/10/2020	A11-MW007 A11-MW007-200910 9/10/2020	A11-MW007** A11-MW007-200910-D 9/10/2020	A11-MW007 A11-MW007-201201 12/2/2020	A11-MW007** A11-MW007-201201-D 12/2/2020
Analyte Name	RG											
Alkalinity, Total (As CaCO3)	--	520	560	540	470	330	370	360	530	530	540	540
Methane	--	20	25	22	14	5.5	3.8	3.9	25 J	21 J	31	29
Nitrate	10	0.2 U	1 U	1 U	0.2 U	0.2 U	0.2 UJ	0.2 UJ	0.12 U	0.12 U	0.12 U	0.12 U
Sulfate	400	10 U	20 U	20 U	13	24	29	28	2.96	2.93	2.45	2.56

Station Location Sample ID Sample Date		A11-MW130A A11-MW130A-200909 9/9/2020	A11-MW130A A11-MW130A-201201 12/1/2020
Analyte Name	RG		
Alkalinity, Total (As CaCO3)	--	330	350
Methane	--	0.00058 UJ	0.0014
Nitrate	10	5.91	6.26
Sulfate	400	17.5 J-	17.3 J-

Notes:  
All results in milligram per liter  
Remediation goals from Record of Decision or Class I Groundwater Standard from 35 IAC 620.410  
Shaded result exceeds remediation goal  
\*\* = Duplicate sample  
U = Not detected above the reported limit  
J = Estimated result J- = Estimated result biased low J+ = Estimated result biased high  
(D) = Diluted result





## Appendix A

# Groundwater Sampling Sheets

## **March 2020 Groundwater Sampling Sheets**



DATE: 3-3-20

WELL #: ~~001~~ Mvood

TIME: 8:28

DEPTH OF PUMP:

WEATHER CONDITIONS: Sunny, cold, 30s

SAMPLERS: Christ T. m

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
0					6.61	1.34	72.2	8.18	11.12	128
5					6.98	1.37	56.4	7.56	11.45	135
10					6.85	1.40	54.0	7.32	11.49	143
15					6.79	1.40	52.7	7.14	11.72	145
20										
25					7.03	1.36	37.4	6.57	12.77	147
30					7.29	1.40	43.9	5.24	11.16	120
35					7.28	1.41	22.6	5.38	11.90	120
40					7.27	1.41	16.7	5.18	11.92	116
45					7.24	1.41	11.9	5.16	11.85	117
50					7.22	1.41	9.3	5.06	12.08	118

\* trouble shooting at down marker well stopped

Sample @ 0120

Fe → 0.24mk



DATE: 3-4-20

WELL #: MW002

TIME: 1255

DEPTH OF PUMP:

WEATHER CONDITIONS:

Sunny, high 40s

SAMPLERS: Tim, Chris

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
0					8.52	0.912	14.7	6.18	13.33	-26
5					7.77	1.12	32.4	1.73	13.45	-157
10					7.77	1.25	61.1	1.60	13.63	-125
15					6.86	1.37	74.3	1.11	13.89	-111
20					6.65	1.37	37.3	1.30	13.63	-120
25					6.62	1.36	21.3	1.04	13.58	-119
30					6.56	1.35	22.9	0.85	13.62	-118
35					6.54	1.36	22.0	0.83	13.61	-116

Sample @ 1335

Fe → 1.76mg/L



SITE NAME: Southeast Rockford, Area 11DATE: 3-4-20

WELL #:

MW003

TIME:

0815 830

DEPTH OF PUMP:

WEATHER CONDITIONS:

Cold, sunny, 40s

SAMPLERS:

Chris + Tim

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
<u>0</u>					<u>6.11</u>	<u>1.06</u>	<u>966</u>	<u>9.09</u>	<u>10.21</u>	<u>-26</u>
<u>0</u>					<u>6.53</u>	<u>1.29</u>	<u>265</u>	<u>4.08</u>	<u>12.17</u>	<u>-89</u>
<u>5</u>					<u>6.58</u>	<u>1.31</u>	<u>177</u>	<u>4.57</u>	<u>11.30</u>	<u>-103</u>
<u>10</u>					<u>6.58</u>	<u>1.32</u>	<u>135</u>	<u>4.60</u>	<u>9.98</u>	<u>-107</u>
<u>15</u>					<u>6.34</u>	<u>1.32</u>	<u>100</u>	<u>3.88</u>	<u>13.74</u>	<u>-99</u>
<u>20</u>					<u>6.63</u>	<u>1.84</u>	<u>23</u>	<u>3.65</u>	<u>12.82</u>	<u>-120</u>
<u>25</u>					<u>6.62</u>	<u>1.84</u>	<u>20.1</u>	<u>3.49</u>	<u>12.43</u>	<u>-120</u>
<u>30</u>					<u>6.62</u>	<u>1.35</u>	<u>17.7</u>	<u>3.39</u>	<u>12.12</u>	<u>-122</u>
<u>35</u>										

Sample @ 905

Fe → 1.32 mL







SITE NAME: Southeast Rockford, Area 11

Sample size 1605

Fe → 0.10 mL

DATE: 3-3-20

TIME: ~~15:10~~ 15:10

WELL #: M0004B

DEPTH OF PUMP:

WEATHER CONDITIONS: Sunny, 40s

SAMPLERS: Chris + Tim

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	pH (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
0					7.57	0.609	0	5.69	12.76	28
08					7.55	0.915	632	1.12	12.83	-1
105					7.73	0.262	315	0.79	12.37	-18
1510					7.53	0.178	262	0.67	12.63	-23
2015					7.37	0.159	259	0.62	12.23	-8
2820										
3015										
3330					6.95	0.233	355	1.00	13.19	12
35					7.05	0.198	428	1.19	12.72	1
40					7.32	1.07	290	1.20	12.39	12
45					7.36	1.14	187	1.46	12.22	29
50					7.42	1.18	100	1.66	12.26	37

\* initially water flowing through the horizon came out very turbid then suddenly (based off appearance) stabilized to 0.0 NTU

\* 1<sup>st</sup> reading is a wash

\* trouble shooting @ 20 ± 25 minute started pump over due to issues



DATE: 3-3-20

WELL #: MWOOS

TIME: 1240

DEPTH OF PUMP:

WEATHER CONDITIONS: Overcast, 40's

SAMPLERS: Chris + Tim

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
0					7.69	0.756	250	5.01	10.89	-72
5					7.71	0.602	196	4.11	<del>10.45</del> 10.45	-66
10					7.51	0.998	65.3	4.28	10.96	-44
15					7.49	1.15	31.3	4.27	10.90	-11
20					7.48	1.21	59.6	4.12	10.96	19
25					7.56	0.846	72.3	4.79	9.95	21
30										
35					7.50	1.14	32.9	5.15	9.92	52
40					7.39	1.27	19.5	4.61	10.47	56
45					7.42	1.29	17.8	4.53	10.62	66
50					7.47	1.27	19.1	4.59	10.99	68
55					7.47	1.25	20.2	4.60	10.36	70

\* @ 30 min mark - Cleaned out horizon due to inconsistencies of turbidity, Saw there was sand & gunk inside the horizon affecting measurements.

Sample @ 1340 Fe → 2002 mg/L



SITE NAME: Southeast Rockford, Area 11

DATE: 3-3-20

WELL #: MW006

TIME: 1030

DEPTH OF PUMP:

WEATHER CONDITIONS: Sunny, 30s

SAMPLERS: Chris & Tim

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
0					7.52	1.34	6.2	7.34	12.11	-108
5					7.55	1.38	5.3	1.72	12.55	-93
10					7.36	1.36	1.8	1.08	12.95	-100
15					7.44	1.36	1.0	0.93	12.93	-110
20					7.43	1.36	0.8	0.85	13.00	-111
25					7.38	1.35	0.8	0.78	12.91	-110

Sample 110:11 Fe → 1.72 mg/L



DATE: 3/4/20  
TIME: 1045

WELL #: MW-007

DEPTH OF PUMP:

WEATHER CONDITIONS: Sunny, 30's

SAMPLERS: Andrew, Chris

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
0					7.03	6.824	78.5	30.84	10.04	-110
5					6.96	6.837	55.7	7.58	9.74	-115
10					6.82	6.858	21.3	6.16	10.12	-120
15					6.81	6.844	14.5	8.02	10.05	-122
20					6.78	6.857	12.7	7.32	10.15	-125
25					6.76	6.863	11.6	6.85	10.11	-126
30					6.76	6.868	9.9	6.74	10.30	-128
35					6.75	6.868	10.1	6.75	10.30	-131
40										

Sample @ 1120

Fe ⇒ 1.80 mg/L



## **June 2020 Groundwater Sampling Sheets**



SITE NAME: Southeast Rockford, Area 11

Fe<sup>2+</sup> 0.10

DATE: 6/19/2020

WELL #: MW-1

TIME: 7:30

DEPTH OF PUMP:

WEATHER CONDITIONS: 75°F, pt. cloudy, 29 mph SAMPLERS: Grobbs + Garbino<sup>av</sup>

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
740			450							
745					6.68	1.20	386	5.75	17.36	56
750					6.70	1.22	250	6.25	17.88	64
755					6.73	1.20	211	5.88	16.50	73
800					6.74	1.22	320	5.98	16.79	78
805				0	6.75	1.21	248	5.52	17.50	80
810					6.74	1.21	173	5.92	17.04	69
815					6.74	1.22	184	5.85	17.22	76
820					6.76	1.21	192	10.90	16.54	67
825					6.74	1.21	115	9.38	17.51	82
830					6.76	1.21	115	8.68	17.58	79
835					6.75	1.21	58.0	8.48	17.73	71

--Difficulty maintaining constant flow rate. Will try different pump on next well.

sample time 08:40



WL - 23.78

Fe<sub>2</sub> 41.58 mg/K (diluted x 2)

WELL #: MW 2

**DEPTH OF PUMP:**

SAMPLERS: Crabs + Carabids

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
11:30			250		6.75	1.35	12.5	3.38	20.30	-101
11:35					6.73	1.35	14.0	2.96	19.54	-102
11:40				0.03	6.71	1.34	13.4	2.81	18.52	-102
11:45					6.67	1.36	22.5	2.74	18.55	-103
11:50					6.67	1.35	22.8	2.62	18.25	-104
11:55					6.67	1.33	22.3	2.48	18.74	-105
12:00					6.65	1.34	20.2	2.54	18.09	-105
12:05					6.65	1.33	19.0	2.50	17.99	-106
12:10	SAMPLES		COLLECTED							



# Well Development Log

**SITE NAME:** Southeast Rockford, Area 11

WL - 24.06

Fe2 - 3.48 mg/l  
(diluted x 2)

DATE: 6/10/2020

TIME: 0700

WELL #: MW 3

**DEPTH OF PUMP:**

WEATHER CONDITIONS: 70°F, cloudy, 5 mph

**SAMPLERS:** Crabs + Caballeros

[illegible]



**SITE NAME:** Southeast Rockford, Area 11

WELL #: NW 4A

**DEPTH OF PUMP:**

WEATHER CONDITIONS: 77° F, partly cloudy, WSW 20 mph

[illegible][illegible]

Fe<sub>2</sub> - 2.39 mg/L



WL 25.10

$\text{Fe}_2: 0.18 \text{ mg/L}$

WELL #: MW-4B

**DEPTH OF PUMP:**

**SAMPLERS:** Grabs + Carabine

[illegible]



2520

29.64

Fe<sub>2</sub>: 0.00 mg/L

WELL #: MW-51

**DEPTH OF PUMP:**

, partly cloudy, ESE 19 mph

[illegible]



# Well Development Log

SITE NAME: Southeast Rockford, Area 11

DATE: 6/9/2020

TIME: 10:12

WEATHER CONDITIONS: 84°F, pt. cloudy, ESE 14 mph

WELL #: MW-6

DEPTH OF PUMP: \_\_\_\_\_

SAMPLERS: Grabs + Carbarino

ELAPSED TIME (MIN)	VOLUME PURGED (mL)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
10:20			400	0	6.87	1.51	14.9	3.56	16.48	-112
10:25					6.84	1.51	16.1	3.52	16.67	-114
10:30					6.83	1.50	18.1	3.45	16.24	-115
10:35					6.82	1.50	17.3	3.41	15.95	-116
10:40					6.80	1.50	14.5	3.36	15.92	-115
10:45					6.80	1.50	10.9	3.23	15.88	-115
10:50					6.80	1.50	7.0	3.23	15.61	-115
10:55					6.80	1.49	4.8	3.11	15.71	-115
11:00	COLLECT SAMPLE									

WL 24.76

Fe<sub>2</sub> = 3.84 mg/L  
(diluted x 2)



323.59

Fe<sub>2</sub> 2.39 mg/K

DATE: 6/10/2020

WELL #: MW 7

TIME: 0911

**DEPTH OF PUMP:**

WEATHER CONDITIONS: 74°F, partly cloudy, 5 mph

[illegible]



## **September 2020 Groundwater Sampling Sheets**



# LOW FLOW GROUNDWATER SAMPLING

SITE NAME: Southeast Rockford, Area 11

Fe<sup>2+</sup> = 0.02 mg/L

DATE: 9/9/20 WELL #: MW-1  
 TIME: 15:15 DEPTH OF PUMP: \_\_\_\_\_  
 WEATHER CONDITIONS: cloudy, wind, cool SAMPLERS: Grabs

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
1535			300		6.42	1.26	300	3.91	14.17	78
1540					6.55	1.27	250	4.18	14.21	84
<del>1545</del> 1545					6.71	1.27	207	4.04	14.26	79
1550					6.77	1.28	185	3.86	14.33	76
1555					6.82	1.28	140	3.82	14.42	73
1600					6.87	1.28	130	4.02	14.51	71
1605					6.89	1.29	111	3.92	14.36	64
1610					6.90	1.28	105	3.80	14.46	65
1615					6.90	1.28	88	3.80	14.34	65
1620					6.93	1.28	81	3.84	14.31	64
1625	SAMPLE									

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Fe<sup>2+</sup>: 2.14 mg/L

## LOW FLOW GROUNDWATER SAMPLING

WL: 24.69

SITE NAME: Southeast Rockford, Area 11

DATE: 9/10/2020

WELL #: MW 2

TIME: 12:07

DEPTH OF PUMP:

WEATHER CONDITIONS: 57°F, overcast, windy

SAMPLERS: O.I.via + MGA

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE <i>ml/min</i>	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
12:20		24.73	250	-0.04	6.68	1.39	34.3	0.79	15.15	-99
12:25					6.69	1.39	34.6	0.57	15.31	-104
12:30					6.69	1.38	27.7	0.53	15.16	-107
12:35					6.71	1.38	22.5	0.43	15.40	-109
12:40					6.70	1.37	18.9	0.40	15.47	-111
12:45					6.71	1.37	18.1	0.37	15.54	-113
12:50					6.70	1.38	13.7	0.39	15.60	-114
12:55					6.70	1.38	13.7	0.36	15.59	-114
13:00					6.71	1.37	13.3	0.33	15.73	-116
13:05	SAMPLES COLLECTED									

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Fe<sup>2+</sup> = 2.52 mg/L

## LOW FLOW GROUNDWATER SAMPLING

WL = 24.99

SITE NAME: Southeast Rockford, Area 11

DATE: 9/10/2020

WELL #: MW-3

TIME: 7:43

DEPTH OF PUMP:

WEATHER CONDITIONS: 50°F, overcast, windy

SAMPLERS: Oлива + MATH

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE mL/min	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
7:55		25.03	400	0.04	7.01	1.28	23.6	3.84	12.82	-99
8:00					7.02	1.29	51.6	1.17	13.28	-123
8:05					7.09	1.29	49.7	0.83	13.32	-131
8:10					7.11	1.29	42.3	0.66	13.36	-136
8:15					7.09	1.29	27.5	0.58	13.47	-137
8:20					7.09	1.28	22.3	0.52	13.57	-138
8:25					7.07	1.28	16.9	0.49	13.60	-138
8:30					7.09	1.28	14.4	0.49	13.56	-140
8:35					7.08	1.27	12.7	0.47	13.68	-140
8:40					7.03	1.26	9.9	0.44	14.13	-140
8:45		SAMPLED								

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.



Fe<sup>2+</sup> = 1.12 mg/L

LOW FLOW GROUNDWATER SAMPLING

WL = 26.37

SITE NAME: Southeast Rockford, Area 11

DATE: 2/10/2020

WELL #: MW-4A

TIME: 14:25

DEPTH OF PUMP:

WEATHER CONDITIONS: 59°F, ~~clear~~ overcast, rain

SAMPLERS: Olivia & Mark

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE ml/min	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
1445		26.38	325	0.01	7.09	1.23	73.1	8.17	14.22	-221
1450					7.12	1.23	52.8	6.85	14.16	-232
1455					7.14	1.23	46.8	6.68	14.11	-236
1500					7.13	1.22	43.6	6.26	14.19	-234
1505					7.14	1.22	39.1	6.08	14.20	-233
1510					7.14	1.23	30.6	5.77	14.21	-231
1515					7.14	1.24	28.6	5.95	14.23	-229
1520					7.14	1.25	22.8	5.80	14.24	-228
1525					7.13	1.26	18.6	5.51	14.25	-225
1530					7.14	1.27	15.1	5.20	14.26	-226
1535					7.13	1.28	13.8	4.91	14.25	-226
1540					7.13	1.29	11.8	4.65	14.26	-226

1545  
Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

1550 SAMPLES COLLECTED

Fe<sup>2+</sup> : less than minimum range      LOW FLOW GROUNDWATER SAMPLING      WL: 26.00

SITE NAME: Southeast Rockford, Area 11

DATE: 9/9/2020

WELL #: NW-4B

TIME: 15:13

DEPTH OF PUMP:

WEATHER CONDITIONS: 57°F, cloudy, windy

SAMPLERS: Matt, Olive

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE mL/min	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
15:30		26.00	300	0.02	7.26	0.304	18.6	3.47	14.40	-11
15:35					6.96	1.04	17.2	2.46	14.17	-61
15:40					7.00	1.21	16.5	2.52	14.00	-24
15:45					6.92	1.15	15.1	2.37	13.91	-26
15:50					6.88	1.21	14.9	2.58	13.86	-12
15:55					6.97	1.24	14.0	2.56	13.93	-4
16:00					7.02	1.26	15.7	2.57	13.97	5
16:05					7.04	1.27	15.7	2.59	14.02	15
16:10					7.06	1.23	14.9	2.70	14.03	25
16:15					7.04	1.24	14.2	2.65	14.00	31
16:20					7.04	1.25	13.5	2.76	14.06	38
16:25					7.04	1.25	13.1	2.80	14.05	41

16:30 SAMPLES COLLECTED

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

# LOW FLOW GROUNDWATER SAMPLING

Fe 2+ : less than minimum range

WL: 25.56

SITE NAME: Southeast Rockford, Area 11

DATE:

9/2/20

TIME:

10:30

WELL #: MW-5

DEPTH OF PUMP:

WEATHER CONDITIONS:

57°F, cloudy, windy

SAMPLERS:

John, Oliver, Matt

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mv)
12:45		25.56	425	0.00	7.07	1.08	200	5.03	15.05	66
12:50					7.04	1.23	165	4.91	14.64	67
12:55					7.01	1.23	137	4.79	14.81	68
13:00					6.90	1.22	67.2	4.75	14.79	78
13:05					6.79	1.23	59.3	4.52	14.47	83
13:10					6.85	1.25	48.4	4.40	14.45	83
13:15					6.75	1.26	43.9	4.10	14.35	85
13:20					6.91	1.25	33.6	3.95	14.26	76
13:25					6.98	1.25	21.8	3.92	14.29	76
13:30					6.99	1.26	16.9	3.83	14.38	77
13:35					7.00	1.26	10.0	3.73	14.43	78
13:40			SAMPLED							

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.



Fe<sup>2+</sup> = 2.19 mg/L

LOW FLOW GROUNDWATER SAMPLING

SITE NAME: Southeast Rockford, Area 11

WL 35, 75

DATE: 9/9/20

WELL #: MW-6

TIME: 11:00

DEPTH OF PUMP:

WEATHER CONDITIONS: 57°F, cloudy

SAMPLERS: Olivia Baker, Matt Carbarney, John Gibbs

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE (mL/min)	DRAWDOWN FEET (+/- 0.3 FT)	pH (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
11:10		25.77	275	0.02	6.92	1.39	21.7	2.39	14.75	-117
11:15					6.91	1.38	17.7	0.86	14.95	-123
11:20					6.91	1.37	16.2	0.67	14.87	-125
11:25					6.95	1.36	17.1	0.56	14.94	-128
11:30					6.95	1.36	13.0	0.54	15.01	-124
11:35					6.82	1.36	13.8	0.54	14.90	-123
11:40					6.94	1.36	12.3	0.49	14.88	-133
11:45	SAMPLES COLLECTED									

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Fe 2+ = over range  
 reduced to 10 mL + filtered to 25 mL distilled (2.5 x)  
 LOW FLOW GROUNDWATER SAMPLING  
 After this reading was 1.99  $2 \times 2.5 = 4.98$

WL = 24.54

DATE: 9/10/2020

WELL #: MW 7

TIME: 9:54

DEPTH OF PUMP:

WEATHER CONDITIONS: 53°F, overcast, windy

SAMPLERS: Olivia & Matt

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
10:05		24.54	400	0.00	6.67	2.22	66.0	2.62	14.34	-97
10:10					6.68	2.13	72.8	0.78	14.43	-107
10:15					6.64	2.05	46.2	0.53	14.51	-113
10:20					6.64	2.01	37.1	0.48	14.73	-117
10:25					6.63	2.01	30.3	0.46	14.77	-120
10:30					6.61	1.98	23.6	0.43	14.93	-121
10:35					6.61	1.96	18.1	0.41	14.93	-124
10:40					6.61	1.94	13.6	0.38	14.85	-125
10:45					6.55	1.94	13.6	0.37	14.88	-123
10:50					6.57	1.94	9.9	0.34	14.82	-126
10:55					SAMPLED					

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

# LOW FLOW GROUNDWATER SAMPLING

SITE NAME: Southeast Rockford, Area 11

WL: 19.18

DATE: 9/9/20

WELL #: MW-130A

TIME: 0830

DEPTH OF PUMP:

WEATHER CONDITIONS: SSF, cloudy, windy  
SAMPLERS: John Gibbs, Matt Gardening, Olivia Burke

Start at 9:00

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
9:05	—	19.56	425	—	—	—	—	—	—	—
9:10			425	0.38	6.68	1.18	89.0	2.46	13.85	17
9:15		19.69		0.51	6.77	1.18	55.9	2.41	13.87	16
9:20					6.82	1.18	36.4	2.39	13.85	18
9:25		19.53			6.88	1.18	25.6	2.33	13.89	20
9:30					6.90	1.18	19.6	2.36	14.14	17
9:35					6.89	1.18	16.0	2.35	14.22	21
9:40					6.77	1.18	15.3	2.42	13.99	24
9:45					6.74	1.18	11.4	2.46	13.91	23
9:50					6.69	1.18	9.8	2.44	13.85	24
9:55	SAMPLES		COLLECTED							

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.



## **December 2020 Groundwater Sampling Sheets**

*Robertson*

LOW FLOW GROUNDWATER SAMPLING

Fe<sup>2+</sup>: 0.19 mg/L  
Static WL 20.16

SITE NAME: Southeast Rockford, Area 11

DATE: 12/01/20  
TIME: 0755  
WELL #: 130A  
DEPTH OF PUMP:  
WEATHER CONDITIONS: clear 19°F  
SAMPLERS: C-60x - A Phelps

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE mL/min	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
8:35	<del>0.1</del>	<del>20.16</del>	350	0.72	7.21	1.10	354	3.75	9.92	78
8:40					7.33	1.14	254	2.15	9.97	38
8:45					7.50	1.16	103	2.60	10.31	-1
8:50					7.61	1.16	131	2.58	10.45	-13
8:55					7.69	1.17	108	2.40	10.57	-20
9:00					7.73	1.17	91.1	2.16	10.57	-23
9:05					7.79	1.17	<del>46.6</del> 79	3.61	10.67	-24
9:10					7.81	1.17	56.4	2.24	10.45	-23
9:15					7.84	1.17	48.7	2.13	10.32	-23
9:20					7.86	1.17	42.7	2.13	10.31	-23
9:25					7.88	1.17	41.1	2.19	10.39	-22
9:30					7.89	1.17	31.8	2.18	10.64	-22

9:35 SAMPLES COLLECTED  
Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

12/01/20 1350  
clear, 340

MW-1

State WL  
28.13

Time	Vol purged	Draw down	pH	Spec cond	Turbidity	DO	temp	ORD
14:01	0.25 gal	475 ml/min	7.96	1.23	537	6.95	12.57	
14:06	1 gal		7.99	1.23	408	5.59	12.84	
14:11	1.5 gal		8.01	1.24	252	5.88	12.84	
14:16	2 gal		8.01	1.23	208	5.84	12.79	
14:21	2.5 gal		8.01	1.23	164	5.79	12.81	
14:26	3.25 gal		8.02	1.23	133	5.72	12.75	
14:31	3.75 gal		8.02	1.23	126	5.61	12.65	
14:36	4 gal + 4.5 gal		8.01	1.23	98.8	6.27	12.75	
14:41	5 gal		8.02	1.22	68.2	5.72	12.64	
14:46	5.25 gal		8.01	1.22	58.0	5.73	12.85	
14:51	5.75 gal		8.01	1.22	49.5	5.81	12.75	
14:56	7 gal		8.01	1.22	36.0	5.78	12.64	
15:01	7.15 gal		8.01	1.22	31.5	5.77	12.65	
15:06	8 gal		8.02	1.20	25.6	6.48	12.40	

pg 1 of 2

COMPUTED BY  
DATE  
PAGE NO.

JOB NO.  
DATE CHECKED  
CHECKED BY

CLIENT  
PROJECT  
DETAIL

CDM  
Smith



PG 2 of 2

WELL #: MW

TIME: \_\_\_\_\_

DEPTH OF PUMP: \_\_\_\_\_

### **SAMPLERS:**

[illegible]

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Ferrus Ion 0.0

# LOW FLOW GROUNDWATER SAMPLING

Fe<sup>2+</sup> = 2.70 mg/L

SITE NAME: Southeast Rockford, Area 11

DATE: 12/2/2020

WELL #: MW002

TIME: 1:58

DEPTH OF PUMP:

WEATHER CONDITIONS: Partly cloudy, 43°F

SAMPLERS: M6H + Alan

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
12:10		25.58	300	0.1	7.14	1.52	125	1.06	13.22	-84
12:15					7.26	1.50	92.0	0.91	13.61	-88
12:20					7.17	1.48	55.4	0.56	13.67	-100
12:25					7.04	1.45	47.7	0.50	13.77	-93
12:30					6.94	1.45	49.8	0.47	13.85	-94
12:35					6.99	1.45	31.8	0.45	13.92	-92
12:40					7.06	1.45	23.5	0.44	13.91	-101
12:45					7.07	1.43	32.7	0.43	14.02	-108
12:50					7.18	1.45	45.1	0.42	14.15	-112
12:55					7.13	1.44	66.2	0.40	14.15	-106
13:00					7.19	1.43	120	0.39	14.14	-111
13:05			350		7.21	1.43	199	0.39	14.18	-109

13:10 SAMPLES COLLECTED

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

# LOW FLOW GROUNDWATER SAMPLING

SITE NAME: Southeast Rockford, Area 11

DATE: 12/2/2020

TIME: 7:17

WELL #: MW 3

DEPTH OF PUMP:

WEATHER CONDITIONS: Partly cloudy 21°F

SAMPLERS: Matt + Alan

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 µS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
800		25.19	375-	-0.67	7.25-	1.35	226	2.19	10.97	-132
<del>830</del> 830					7.42	1.35	104	1.16	10.70	-143
810					7.38	1.35	55-1	0.91	11.23	-142
815					7.31	1.35	32.4	0.82	11.31	-141
820					7.41	1.35	20.8	0.87	10.64	-144
825					7.49	1.35	14.4	0.68	11.3	-154
830					7.56	1.35	11.5	0.68	11.17	-155
835					7.49	1.34	9.4	0.67	11.83	-157
840					7.55	1.34	7.0	0.62	11.75	-158
845	SAMPLES COLLECTED									

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Dilution factor  
 $Fe^{2+} = 1.54 \times 2.5 = 3.85 \text{ mg/L}$



# LOW FLOW GROUNDWATER SAMPLING

SITE NAME: Southeast Rockford, Area 11

DATE: 12/2/2020

TIME: 14:04

WELL #: MW 41A

DEPTH OF PUMP:

WEATHER CONDITIONS: Partly Cloudy 44°F

SAMPLERS: Matt + Alan

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
1420		27.27	330	0.39	7.59	1.31	155	5.52	13.98	-292
1425					7.75	1.30	65.3	4.11	13.65	-298
1430					7.53	1.30	35.2	3.29	13.41	-284
1435					7.44	1.31	17.7	2.85	13.17	-280
1440					7.73	1.31	9.2	0.54	12.35	-290
1445					7.58	1.32	7.5	0.58	12.73	-276
1450					7.59	1.33	5.8	0.56	12.56	-279
1455					7.43	1.34	3.6	0.54	12.34	-276
1500	SAMPLES COLLECTED									

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

# LOW FLOW GROUNDWATER SAMPLING

SITE NAME: Southeast Rockford, Area 11

DATE: 12/1/2020

TIME: 15:12

WELL #: MW 4B

DEPTH OF PUMP:

WEATHER CONDITIONS: clear 34°F

SAMPLERS: Matt Carlen

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 µS/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
15:40		<del>26.95</del> 26.95	400	0.05	7.19	1.29	964	16.61	9.46	41
15:45					7.29	1.27	506	8.46	9.37	50
15:50					7.31	1.32	159	7.40	9.93	56
15:55					7.35	1.38	114	6.69	9.95	59
16:00					7.43	1.40	99.8	6.28	9.39	58
16:05					7.41	1.40	97.2	5.60	8.75	67
16:10					7.39	1.40	108	5.13	9.90	65
16:20					7.38	1.41	74.0	4.97	10.17	82
16:25					7.38	1.41	52.0	4.74	10.32	98
16:30					7.37	1.39	40.8	4.51	10.72	95
16:35					7.36	1.39	31.1	4.36	10.74	99
16:40					7.39	1.40	32.0	4.15	10.93	94

16:45 SAMPLES COLLECTED

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Fe<sup>2+</sup>: 0.73 mg/L

# LOW FLOW GROUNDWATER SAMPLING

Fe<sup>2+</sup>: 0.03 mg/L

SITE NAME: Southeast Rockford, Area 11

DATE: 12/1/2020

WELL #: MW 005

TIME: 12:41

DEPTH OF PUMP:

WEATHER CONDITIONS: Clear 32°F

SAMPLERS: Matt Carbanas

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 mS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 C°)	REDOX POTENTIAL mV (+/- 10 mV)
1250		26.47	500	0.2	6.40	1.41	237	9.00	16.12	1285
1255					6.92	1.49	313	5.62	11.51	137
1300					7.01	1.48	267	5.56	11.49	141
1305					7.16	1.48	199	5.27	10.18	146
1310					7.13	1.48	126	5.17	11.48	150
1315					7.15	1.47	106	5.08	11.83	152
1320					7.15	1.47	85.7	5.08	11.62	155
1325					7.15	1.48	68.5	5.02	11.45	157
1330					7.16	1.49	64.8	5.12	10.93	156
1335					7.14	1.50	45.6	5.03	12.15	159
1340					7.15	1.50	41.1	4.96	12.01	159
1345					7.16	1.49	39.9	4.85	12.84	160

1350 SAMPLES COLLECTED

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.



# LOW FLOW GROUNDWATER SAMPLING

Stable water level  
26.57

SITE NAME: Southeast Rockford, Area 11

DATE: 12/01

WELL #: NW -6

TIME: 1105

DEPTH OF PUMP:

WEATHER CONDITIONS: clear 30s

SAMPLERS:

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 $\mu$ S/cm)	TURBIDITY NTUS (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP $^{\circ}$ C (+/- 5 $^{\circ}$ C)	REDOX POTENTIAL mV (+/- 10 mV)
1126	1 gal		500 ml/min		7.87	1.43	76.5	3.04	12.35	-134
1131	2 gal				7.88	1.43	23.1	1.55	12.62	-139
1136	2.5 gal				7.90	1.41	13.6	1.48	12.61	-142
1141	3.5 gal				7.91	1.40	11.5	1.45	12.64	-145
1146	4.5 gal				7.91	1.39	8.4	1.43	12.36	-146
1151	4.75 gal				7.91	1.39	6.4	1.39	12.73	-147
1156	5 gal				8.03	1.39	5.0	1.42	12.62	-149
1201	5.25 gal				7.94	1.38	3.8	1.35	12.35	-150
1206	5.5 gal				7.92	1.38	2.7	1.33	12.66	-150
1211			SAMPLE							

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.

Ferris 266

# LOW FLOW GROUNDWATER SAMPLING

WL 25-41

SITE NAME: Southeast Rockford, Area 11

Fe<sup>2+</sup> = 2.86 mg/L

DATE: 12/2/2020

WELL #: MW 007

TIME: 1957

DEPTH OF PUMP:

WEATHER CONDITIONS: Clear 33°F

SAMPLERS: Alan + Matt

ELAPSED TIME (MIN)	VOLUME PURGED (GALS)	DEPTH TO WATER (FT TIC)	FLOW RATE	DRAWDOWN FEET (+/- 0.3 FT)	ph (+/- 0.25 SU)	SPECIFIC COND. (+/- 50 µS/cm)	TURBIDITY NTUs (+/- 10%)	DISSOLVED OXYGEN mg/L (+/- 10%)	TEMP °C (+/- 5 °C)	REDOX POTENTIAL mV (+/- 10 mV)
1010		25.41	450	0.03	6.95	1.43	194	6.79	10.80	-81
1015					6.86	1.39	206	3.17	11.65	-92
1020					6.87	1.39	125	2.54	11.84	-106
1025					6.93	1.36	74.5	1.46	11.97	-116
1030					6.97	1.37	44.4	1.15	12.31	-118
1035					6.96	1.37	31.8	0.57	12.16	-126
1040					6.95	1.38	31.3	0.54	12.35	-127
1045					6.86	1.39	28.4	0.53	12.28	-124
1050	SAMPLES COLLECTED				7.02	1.38	20.7	0.50	12.57	-128

Drawdown is not to exceed 0.3 of a foot. Flow rate should not exceed 500 ml/min during purging or 250 ml/min during sampling. Readings should be taken every three to five minutes. The well is considered stabilized and ready for sampling when the indicator parameters have stabilized for three consecutive readings by the measurements indicated in parentheses. Turbidity +/- 10% or less than 5 NTU.



## Appendix B

# Data Validation Reports and Data Packages



## **March 2020 Data Validation Reports and Data Packages**

**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** E200303  
**Laboratory:** ESAT / Tech Law  
**Matrix:** Groundwater  
**Collection date:** 03/03/2020 & 03/04/2020  
**Analysis/Methods:** 1,4-Dioxane - SW-846 8000D SIM

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
E200303-01	A11-TB002-200304	E200303-07	A11-MW004B-200303
E200303-02	A11-FB001-200303	E200303-08	A11-MW005-200303
E200303-03	A11-MW001-200303	E200303-09	A11-MW006-200303
E200303-04	A11-MW002-200304	E200303-10	A11-MW007-200304
E200303-05	A11-MW003-200304	E200303-11	A11-MW007-200304-D
E200303-06	A11-MW004A-200304	E200303-12	A11-TB001-200303

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Volatile Organic Compounds 8260 / 1,4-Dioxane 8000D**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	Yes		
Laboratory Control Spike Duplicates RPD within limits?	No		
Laboratory Duplicate RPDs within limits?	N/A		
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b><u>Sample</u> A11-MW007-200304</b>	<b><u>Duplicate</u> A11-MW007-200304-D</b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
			Acceptable		

<b>MS/MSD</b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
E20C007-MS1 / MSD1 (200303-04)	Acceptable			
E20C007-MS2 / MSD2 (200303-08)	Acceptable			

<b>LCS/LCSD</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	No		
Laboratory Control Sample criteria met?	No		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	No		
Was the ICAL criteria met?	No		
Was the CCV criteria met?	No		
Was the Tuning criteria met?	Yes		
Were the Surrogate % recoveries within laboratory determined control limits?	Yes		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

<b>Blanks</b>	<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
E20C007-BLK1	Nondetect			

<b>Field Blank</b> A11-FB001-200303 A11-TB001-200303 A11-TB002-200304		<u>Concentration</u> Nondetect Nondetect Nondetect	<u>MDL / PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>Surrogates</b>		<u>%R</u> Acceptable	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>MS/MSD</b> E20C007-MS1 / MSD1 (200303-04) E20C007-MS2 / MSD2 (200303-08)		<u>%R</u> Acceptable  Acceptable	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>LCS/LCSD</b> E20C007-BS1		<u>%R</u> Acceptable	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>ICAL</b> November 20, 2019		<u>RRF</u> Acceptable	<u>%RSD</u> Acceptable	<u>Limits</u>	<u>Qualifiers</u> <u>Associated Samples</u>
<b>ICV / CCV</b> 3/10/2020 10:44 3/10/2020 3:41 3/11/2020 10:14 3/11/2020 1:23		<u>RRF</u> Acceptable Acceptable Acceptable Acceptable	<u>%D</u> Acceptable Acceptable Acceptable Acceptable	<u>Limits</u>	<u>Qualifiers</u> <u>Associated Samples</u>
<b>Tune</b> Acceptable					
<b>MRL Check</b>  E20C007-MRL1		<u>%R</u>  Acceptable	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>Internal Standards</b>		<u>Area</u>	<u>Area Lower / Upper</u> <u>Limit</u> Acceptable	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>Representativeness:</b>					
Were sampling procedures and design criteria met?					<u>Yes</u> <u>No</u> <u>N/A</u>
Were holding times met?					Yes
Was preservation criteria met? (0° C - 6° C)					Yes
Were Chain-of-Custody records complete and provided in data package?					Yes
<u>Comments (note deviations):</u> The cooler temperature was -0.8 ° C.					
<b>Preservation</b>		<u>Cooler Temperature (Degrees C)</u> Acceptable	<u>Preservation Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
<b>Holding Times</b>	<u>Analyte</u>	<u>Days to Extraction</u> Acceptable	<u>HT Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
<b>Comparability:</b>					
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?					<u>Yes</u> <u>No</u> <u>N/A</u>
<u>Comments (note deviations):</u>					Yes
<b>Completeness (90%):</b>					
Are all data in this SDG usable?					<u>Yes</u> <u>No</u> <u>N/A</u>
<u>Comments (note deviations):</u>					Yes

**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):**Yes No N/A****Yes****Yes****Comment:**

Data is usable as reported.

Data Validator:

*Kristine Molloy*Date: *4/3/2020*

Data Reviewer:

Cherie ZakowskiDate: 4/5/2020



Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Howard Pham

**Reported:**  
Mar-18-20 15:26

## 1,4-Dioxane by GC-MS

### TechLaw - ESAT Contract

#### A11-TB002-200304 (E200303-01)

Matrix: Water

Sampled: Mar-04-20 08:00

Received: Mar-05-20 13:30

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.221	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.780			70.5%		70-130	"	"	"

#### A11-FB001-200303 (E200303-02)

Matrix: Water

Sampled: Mar-03-20 17:00

Received: Mar-05-20 13:30

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.202	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.759			75.3%		70-130	"	"	"

#### A11-MW001-200303 (E200303-03)

Matrix: Water

Sampled: Mar-03-20 09:20

Received: Mar-05-20 13:30

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>6.85</b>			0.202	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.802			79.6%		70-130	"	"	"

#### A11-MW002-200304 (E200303-04)

Matrix: Water

Sampled: Mar-04-20 13:35

Received: Mar-05-20 13:30

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>3.31</b>			0.203	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.891			87.7%		70-130	"	"	"

#### A11-MW003-200304 (E200303-05)

Matrix: Water

Sampled: Mar-04-20 09:05

Received: Mar-05-20 13:30

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>8.57</b>			0.202	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.822			81.6%		70-130	"	"	"

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Project Manager: Howard Pham

**Reported:**  
Mar-18-20 15:26

## 1,4-Dioxane by GC-MS

### TechLaw - ESAT Contract

#### A11-MW004A-200304 (E200303-06)

**Matrix: Water**

**Sampled: Mar-04-20 15:20**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>1.41</b>			0.203	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.789			77.6%		70-130	"	"	"

#### A11-MW004B-200303 (E200303-07)

**Matrix: Water**

**Sampled: Mar-03-20 16:05**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>9.75</b>			0.202	ug/L	1	E20C007	Mar-09-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.784			77.8%		70-130	"	"	"

#### A11-MW005-200303 (E200303-08)

**Matrix: Water**

**Sampled: Mar-03-20 13:40**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>5.35</b>			0.202	ug/L	1	E20C007	Mar-09-20	Mar-11-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.856			84.9%		70-130	"	"	"

#### A11-MW006-200303 (E200303-09)

**Matrix: Water**

**Sampled: Mar-03-20 11:00**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>1.54</b>			0.202	ug/L	1	E20C007	Mar-09-20	Mar-11-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.819			81.2%		70-130	"	"	"

#### A11-MW007-200304 (E200303-10)

**Matrix: Water**

**Sampled: Mar-04-20 11:20**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>3.38</b>			0.203	ug/L	1	E20C007	Mar-09-20	Mar-11-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.861			84.7%		70-130	"	"	"

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Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
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 Project Manager: Howard Pham

**Reported:**  
 Mar-18-20 15:26

## 1,4-Dioxane by GC-MS

### TechLaw - ESAT Contract

**A11-MW007-200304-D (E200303-11)**

**Matrix: Water**

**Sampled: Mar-04-20 11:20**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>3.30</b>			0.205	ug/L	1	E20C007	Mar-09-20	Mar-11-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.997			97.3%		70-130	"	"	"

**A11-TB001-200303 (E200303-12)**

**Matrix: Water**

**Sampled: Mar-03-20 08:00**

**Received: Mar-05-20 13:30**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.207	ug/L	1	E20C007	Mar-09-20	Mar-11-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.848			82.0%		70-130	"	"	"

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Reported:  
Mar-18-20 15:26

## 1,4-Dioxane by GC-MS - Quality Control

### TechLaw - ESAT Contract

#### Batch E20C007 - EPA 522

##### Blank (E20C007-BLK1)

Prepared: Mar-09-20 Analyzed: Mar-10-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	U			0.200	ug/L						
Surrogate: 1,4-Dioxane-d8	0.766				"	1.00		76.6%	70-130		

##### LCS (E20C007-BS1)

Prepared: Mar-09-20 Analyzed: Mar-10-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	0.764			0.200	ug/L	1.00		76.4%	70-130		
Surrogate: 1,4-Dioxane-d8	0.728				"	1.00		72.8%	70-130		

##### MRL Check (E20C007-MRL1)

Prepared: Mar-09-20 Analyzed: Mar-10-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	U	Q		0.200	ug/L	0.200		%	50-150		
Surrogate: 1,4-Dioxane-d8	0.738				"	1.00		73.8%	70-130		

##### Matrix Spike (E20C007-MS1)

Source: E200303-04

Prepared: Mar-09-20 Analyzed: Mar-10-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	4.19			0.207	ug/L	1.03	3.31	85.7%	70-130		
Surrogate: 1,4-Dioxane-d8	0.911				"	1.03		88.2%	70-130		

##### Matrix Spike (E20C007-MS2)

Source: E200303-08

Prepared: Mar-09-20 Analyzed: Mar-11-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	6.32			0.203	ug/L	1.02	5.35	95.4%	70-130		
Surrogate: 1,4-Dioxane-d8	0.822				"	1.02		80.9%	70-130		

##### Matrix Spike Dup (E20C007-MSD1)

Source: E200303-04

Prepared: Mar-09-20 Analyzed: Mar-10-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	4.16			0.205	ug/L	1.02	3.31	83.4%	70-130	2.66	30
Surrogate: 1,4-Dioxane-d8	0.880				"	1.02		85.9%	70-130		





TechLaw Inc ESAT Region 5  
536 South Clark Street, Suite 734  
Chicago, IL 60605  
(312) 353-8303  
(312) 353-5814 (Fax)  
[www.techlawinc.com](http://www.techlawinc.com)

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Howard Pham

**Reported:**  
Mar-18-20 15:26

**1,4-Dioxane by GC-MS - Quality Control**  
**TechLaw - ESAT Contract**

**Batch E20C007 - EPA 522**

Matrix Spike Dup (E20C007-MSD2)

Source: E200303-08

Prepared: Mar-09-20 Analyzed: Mar-11-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	6.34			0.203	ug/L	1.02	5.35	97.0%	70-130	1.71	30
Surrogate: 1,4-Dioxane-d8	0.859				"	1.02		84.6%	70-130		

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Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Howard Pham

**Reported:**  
Mar-18-20 15:26

## Notes and Definitions

U Not Detected  
NR Not Reported  
Q QC limit Exceeded

**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** E200306  
**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch  
**Matrix:** Groundwater  
**Collection date:** 03/03/2020 & 03/04/2020  
**Analysis/Methods:** Volatile Organic Compounds (VOCs) 8260

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
2003006-01	A11-TB002-200304	2003006-07	A11-MW004B-200303
2003006-02	A11-FB001-200303	2003006-08	A11-MW005-200303
2003006-03	A11-MW001-200303	2003006-09	A11-MW006-200303
2003006-04	A11-MW002-200304	2003006-10	A11-MW007-200304
2003006-05	A11-MW003-200304	2003006-11	A11-MW007-200304-D
2003006-06	A11-MW004A-200304	2003006-12	A11-TB001-200303

Data validation was performed in accordance with the specific analytical method and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Volatile Organic Compounds 8260 / 1,4-Dioxane 8000D**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?		No	
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)		Yes	
Laboratory Control Spike Duplicates RPD within limits?		No	
Laboratory Duplicate RPDs within limits?		N/A	
<u>Comments (note deviations):</u>			

Field Duplicates	Sample A11-MW007-200304	Duplicate A11-MW007-200304-D	%RPD	Qualifiers	Associated Samples
1,3,5-Trimethylbenzene	4.84	4.57	NC	None	Sample results < 5xs RL; ABS Diff. < RL
Isopropylbenzene	12	11	NC	None	
n-Propylbenzene	6.68	6.03	NC	None	

MS/MSD	%RPD	Limit	Qualifiers	Associated Samples
B20C019-MS1 / MSD1 (2003006-04RE1)	Acceptable			
B20D016-MS1 / MSD1 (2003006-08RE2)	Acceptable			

LCS/LCSD	%RPD	Limits	Qualifiers	Associated Samples
B20C012-BS1 / BSD1	Acceptable			
B20C019-BS1 / BSD1	1,1-Dichloroethene 2,2-Dichloropropane Hexachlorobutadiene n-Butylbenzene	21.8 55.3 21.3 23.2	20% 20% 20% 20%	J** J** J** J** 2003006-03RE1 through 2003006-06RE1, 2003006-10RE1, 2003006-11RE1
B20C016-BS1	Acceptable			

\*\*Sample results nondetect - no qualifiers required.

Laboratory Duplicate	%RPD	Limits	Qualifiers	Associated Samples
N/A				

<b>Accuracy:</b>		<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)			No	
Laboratory Control Sample criteria met?			No	
Were the Laboratory Method Blank results all < RL?			Yes	
Were the Field Blanks results all < RL?			No	
Was the ICAL criteria met?			No	
Was the CCV criteria met?			No	
Was the Tuning criteria met?			Yes	
Were the Surrogate % recoveries within laboratory determined control limits?			Yes	
Were the Internal Standard areas within ± 50 - 150%?			N/A	
<u>Comments (note deviations):</u>				

<b>Blanks</b>	<b>Concentration</b>	<b>MDL /PQL</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
E20C012-BLK1	Nondetect			
E20C012-BLK2	Nondetect			
E20C019-BLK1	Nondetect			
E20C019-BLK2	Nondetect			
E20C016-BLK1	Nondetect			

Field Blank		<u>Concentration</u>	<u>MDL / PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
A11-FB001-200303		Nondetect			
A11-TB001-200303	cis-1,2-Dichloroethene	2.33	0.43 / 2.0	None	Sample results nondetect
A11-TB002-200304	cis-1,2-Dichloroethene	2.2	0.43 / 2.0	None	Sample results nondetect

<b>Surrogates</b>	<b>%R</b>	<b>Limit</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
	Acceptable			

MS/MSD		%R	Limits (%)	Qualifiers	Associated Samples
B20C019-MS1 / MSD1 (2003006-04RE1)	Carbon Disulfide	60.5 / 58.6	60-110	J / UJ	2003006-04RE1**
B20D016-MS1 / MSD1 (2003006-08RE2)		Acceptable			

\*\*Results reported from 2003006-04 - no qualification required

LCS/LCSD		%R	Limits	Qualifiers	Associated Samples
B20C012-BS1 / BSD1		Acceptable			
B20C019-BS1 / BSD1					
	Acetone	146 / 136	70-130	J**	2003006-03RE1 through 2003006-06RE1, 2003006-10RE1, 2003006-11RE1
	2,2-Dichloropropane	110 / 62.6	70-130	J / UJ	
B20D016-BS1	Bromomethane	65.2	70-130	J / UJ	2003006-05RE2 through 2003006-09RE2

\*\*Sample results nondetect - no qualifiers required.

ICAL		RRF	%RSD	Limits	Qualifiers	Associated Samples
3/5/2020 13:53	1,1-Dichloroethene	Acceptable	22.52	20	J**	All samples
	Carbon Disulfide	Acceptable	20.18	20	J**	All samples
	Carbon Tetrachloride	Acceptable	26.45	20	J**	All samples
	Tetrachloroethene	Acceptable	21.62	20	J**	All samples
	1,1,1-Trichloroethane	Acceptable	20.67	20	J	All samples

\*\*Sample results nondetect - no qualifiers required.

ICV / CCV			RRF	%D	Limits	Qualifiers	Associated Samples
ICV							
3/05/2020	17:56	Acetone	Acceptable	96.5	40	J / UJ	All samples
3/05/2020 2:38			Acceptable	Acceptable			



CCV						
3/09/2020 8:39		Acceptable	Acceptable			
3/09/2020 17:15		Acceptable	Acceptable			
3/10/2020 17:12	Acetone	Acceptable	-46.5	40	J / UJ	2003006-03RE1 through 2003006-06RE1, 2003006-10RE1, 2003006-11RE1
3/11/2020 4:24	Acetone	Acceptable	-35.6	40	J / UJ	
3/11/2020 10:40	Bromomethane	Acceptable	35.2	30	J / UJ	2003006-05RE2 through 2003006-09RE2
	Carbon Disulfide	Acceptable	28	25	J / UJ	
	Trans 1,3-Dichloropropane	Acceptable	20.2	20	J / UJ	
<b>MRL Check</b>			<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
B20C012-MRL1			Acceptable			
<b>Tune</b>						
Acceptable						
<b>Internal Standards</b>						
	<u>Area</u>	<u>Area Lower / Upper Limit</u>			<u>Qualifiers</u>	<u>Associated Samples</u>
		Acceptable				
<b>Representativeness:</b>						<u>Yes No N/A</u>
Were sampling procedures and design criteria met?						Yes
Were holding times met?						Yes
Was preservation criteria met? (0° C - 6° C)						Yes
Were Chain-of-Custody records complete and provided in data package?						Yes
<u>Comments (note deviations):</u> The cooler temperatures were 4.7 & 5.6 ° C.						
<b>Preservation</b>	<u>Cooler Temperature (Degrees C)</u>	<u>Preservation Criteria</u>		<u>Qualifier</u>	<u>Associated Samples</u>	
	Acceptable					
<b>Holding Times</b>	<u>Analyte</u>	<u>Days to Extraction</u>	<u>HT Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>	
		Acceptable				
<b>Comparability:</b>						<u>Yes No N/A</u>
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?						Yes
<u>Comments (note deviations):</u>						
<b>Completeness (90%):</b>						<u>Yes No N/A</u>
Are all data in this SDG usable?						Yes
<u>Comments (note deviations):</u>						
<b>Sensitivity:</b>						<u>Yes No N/A</u>
Are MDLs present and reported?						Yes
Do the reporting limits meet project requirements?						Yes
<u>Comments (note deviations):</u>						
<b>Comment:</b>						
As stated in the case narrative, samples were first screened at a 50x dilution and several dilution factors were subsequently required. Each analyte is reported at the lowest dilution factor for which the analyte concentration remained within calibration range.						
Case narrative indicates co-elution affected the calculated concentration of n-butylbenzene in sample 2003006-05RE2. Potential bias of the quantification ion led to the result being qualified as estimated by the laboratory.						
As stated in the case narrative, no -BSD1 is associated with the B20D016 batch due to an error in laboratory instrument that was not noticed or corrected until after the time frame required by the SOP for a closing CCV.						
Data is usable with appropriate qualifiers applied.						
Data Validator:	<u>Kristine Molloy</u>			Date:	<u>4/16/2020</u>	
Data Reviewer:	<u>Cherie Zakowski</u>			Date:	<u>4/18/2020</u>	



**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-TB002-200304 (2003006-01)**

**Matrix: Water**

**Sampled: Mar-04-20 08:00**

**Received: Mar-05-20 14:45**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	2.20			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"



Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB002-200304 (2003006-01)

Matrix: Water

Sampled: Mar-04-20 08:00

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Chlorobenzene	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	11.7			117%		73-124	"	"	"
1,2-Dichloroethane-d4	10.9			108%		84-122	"	"	"
Toluene-d8	10.6			106%		88-108	"	"	"
4-Bromofluorobenzene	9.12			91.2%		84-108	"	"	"



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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-FB001-200303 (2003006-02)

Matrix: Water

Sampled: Mar-03-20 17:00

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-FB001-200303 (2003006-02)

Matrix: Water

Sampled: Mar-03-20 17:00

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.8			108%		73-124	"	"	"
1,2-Dichloroethane-d4	10.7			107%		84-122	"	"	"
Toluene-d8	10.0			100%		88-108	"	"	"
4-Bromofluorobenzene	9.28			92.8%		84-108	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-MW001-200303 (2003006-03RE1)

Matrix: Water

Sampled: Mar-03-20 09:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20C019	Mar-10-20	Mar-10-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	4.51			2.00	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	6.74			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW001-200303 (2003006-03RE1)

Matrix: Water

Sampled: Mar-03-20 09:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20C019	Mar-10-20	Mar-10-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.3			102%		73-124	"	"	"
1,2-Dichloroethane-d4	11.1			110%		84-122	"	"	"
Toluene-d8	9.78			97.8%		88-108	"	"	"
4-Bromofluorobenzene	9.03			90.3%		84-108	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-MW002-200304 (2003006-04)

Matrix: Water

Sampled: Mar-04-20 13:35

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			100	ug/L	50	B20C012	Mar-06-20	Mar-09-20
Chloromethane	U			100	"	"	"	"	"
Vinyl chloride	U			100	"	"	"	"	"
Bromomethane	U			100	"	"	"	"	"
Chloroethane	U			100	"	"	"	"	"
Trichlorofluoromethane	U			100	"	"	"	"	"
1,1-Dichloroethene	U			100	"	"	"	"	"
Acetone	U			625	"	"	"	"	"
Carbon disulfide	U	(MS), L		100	"	"	"	"	"
Methylene chloride	U			100	"	"	"	"	"
trans-1,2-Dichloroethene	U			100	"	"	"	"	"
1,1-Dichloroethane	U			100	"	"	"	"	"
2,2-Dichloropropane	U			100	"	"	"	"	"
cis-1,2-Dichloroethene	U			100	"	"	"	"	"
2-Butanone	U			625	"	"	"	"	"
Bromochloromethane	U			100	"	"	"	"	"
Chloroform	U			100	"	"	"	"	"
1,1,1-Trichloroethane	U			100	"	"	"	"	"
Carbon tetrachloride	U			100	"	"	"	"	"
1,1-Dichloropropene	U			100	"	"	"	"	"
Benzene	U			100	"	"	"	"	"
1,2-Dichloroethane	U			100	"	"	"	"	"
Trichloroethene	U			100	"	"	"	"	"
1,2-Dichloropropane	U			100	"	"	"	"	"
Dibromomethane	U			100	"	"	"	"	"
Bromodichloromethane	U			100	"	"	"	"	"
cis-1,3-Dichloropropene	U			100	"	"	"	"	"
4-Methyl-2-pentanone	U			250	"	"	"	"	"
trans-1,3-Dichloropropene	U			100	"	"	"	"	"
1,1,2-Trichloroethane	U			100	"	"	"	"	"
Tetrachloroethene	U			100	"	"	"	"	"
1,3-Dichloropropane	U			100	"	"	"	"	"
2-Hexanone	U			250	"	"	"	"	"
Dibromochloromethane	U			100	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			100	"	"	"	"	"
Chlorobenzene	U			100	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			100	"	"	"	"	"





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Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW002-200304 (2003006-04)

Matrix: Water

Sampled: Mar-04-20 13:35

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Styrene	U			100	ug/L	50	B20C012	Mar-06-20	Mar-09-20
Bromoform	U			100	"	"	"	"	"
Isopropylbenzene	121			100	"	"	"	"	"
Bromobenzene	U			100	"	"	"	"	"
1,2,3-Trichloropropane	U			100	"	"	"	"	"
n-Propylbenzene	215			100	"	"	"	"	"
2-Chlorotoluene	U			100	"	"	"	"	"
1,3,5-Trimethylbenzene	285			100	"	"	"	"	"
4-Chlorotoluene	U			100	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			100	"	"	"	"	"
tert-Butylbenzene	U			100	"	"	"	"	"
1,2,4-Trimethylbenzene	822			100	"	"	"	"	"
sec-Butylbenzene	U			100	"	"	"	"	"
1,3-Dichlorobenzene	U			100	"	"	"	"	"
p-Isopropyltoluene	U			100	"	"	"	"	"
1,4-Dichlorobenzene	U			100	"	"	"	"	"
1,2-Dichlorobenzene	U			100	"	"	"	"	"
n-Butylbenzene	U			100	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			100	"	"	"	"	"
1,2,4-Trichlorobenzene	U			100	"	"	"	"	"
Hexachlorobutadiene	U			100	"	"	"	"	"
Naphthalene	U			100	"	"	"	"	"
1,2,3-Trichlorobenzene	U			100	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.5			105%		73-124	"	"	"
1,2-Dichloroethane-d4	10.6			105%		84-122	"	"	"
Toluene-d8	10.3			103%		88-108	"	"	"
4-Bromofluorobenzene	10.6			106%		84-108	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW002-200304 (2003006-04RE1)

Matrix: Water

Sampled: Mar-04-20 13:35

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Toluene</b>	<b>78600</b>			4000	ug/L	2000	B20C019	Mar-10-20	Mar-10-20
<b>Ethylbenzene</b>	<b>6840</b>			4000	"	"	"	"	"
<b>m+p-Xylene</b>	<b>19800</b>			8000	"	"	"	"	"
<b>o-Xylene</b>	<b>5100</b>			4000	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>Dibromofluoromethane</i>	10.3			102%		73-124	"	"	"
<i>1,2-Dichloroethane-d4</i>	11.1			110%		84-122	"	"	"
<i>Toluene-d8</i>	9.76			97.6%		88-108	"	"	"
<i>4-Bromofluorobenzene</i>	9.45			94.5%		84-108	"	"	"

A11-MW003-200304 (2003006-05RE1)

Matrix: Water

Sampled: Mar-04-20 09:05

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Ethylbenzene</b>	<b>1500</b>			400	ug/L	200	B20C019	Mar-10-20	Mar-10-20
<b>m+p-Xylene</b>	<b>13000</b>			800	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>Dibromofluoromethane</i>	10.4			104%		73-124	"	"	"
<i>1,2-Dichloroethane-d4</i>	11.2			111%		84-122	"	"	"
<i>Toluene-d8</i>	9.78			97.8%		88-108	"	"	"
<i>4-Bromofluorobenzene</i>	9.50			95.0%		84-108	"	"	"

A11-MW003-200304 (2003006-05RE2)

Matrix: Water

Sampled: Mar-04-20 09:05

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Dichlorodifluoromethane</b>	U			20.0	ug/L	10	B20D016	Mar-10-20	Mar-11-20
<b>Chloromethane</b>	U			20.0	"	"	"	"	"
<b>Vinyl chloride</b>	U			20.0	"	"	"	"	"
<b>Bromomethane</b>	U	(LCS), J		20.0	"	"	"	"	"
<b>Chloroethane</b>	U			20.0	"	"	"	"	"
<b>Trichlorofluoromethane</b>	U			20.0	"	"	"	"	"
<b>1,1-Dichloroethene</b>	U			20.0	"	"	"	"	"
<b>Acetone</b>	U			125	"	"	"	"	"
<b>Carbon disulfide</b>	U			20.0	"	"	"	"	"





## Environmental Protection Agency Region 5

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536 South Clark Street, Chicago, IL 60605  
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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

### Volatiles by GC/MS, EPA 8260C (modified)

### US EPA Region 5 LSASD Analytical Services Branch

A11-MW003-200304 (2003006-05RE2)

Matrix: Water

Sampled: Mar-04-20 09:05

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Methylene chloride	U			20.0	ug/L	10	B20D016	Mar-10-20	Mar-11-20
trans-1,2-Dichloroethene	U			20.0	"	"	"	"	"
1,1-Dichloroethane	U			20.0	"	"	"	"	"
2,2-Dichloropropane	U			20.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			20.0	"	"	"	"	"
2-Butanone	U			125	"	"	"	"	"
Bromochloromethane	U			20.0	"	"	"	"	"
Chloroform	U			20.0	"	"	"	"	"
1,1,1-Trichloroethane	U			20.0	"	"	"	"	"
Carbon tetrachloride	U			20.0	"	"	"	"	"
1,1-Dichloropropene	U			20.0	"	"	"	"	"
Benzene	U			20.0	"	"	"	"	"
1,2-Dichloroethane	U			20.0	"	"	"	"	"
Trichloroethene	U			20.0	"	"	"	"	"
1,2-Dichloropropane	U			20.0	"	"	"	"	"
Dibromomethane	U			20.0	"	"	"	"	"
Bromodichloromethane	U			20.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			20.0	"	"	"	"	"
4-Methyl-2-pentanone	U			50.0	"	"	"	"	"
Toluene	38.4			20.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			20.0	"	"	"	"	"
1,1,2-Trichloroethane	U			20.0	"	"	"	"	"
Tetrachloroethene	U			20.0	"	"	"	"	"
1,3-Dichloropropane	U			20.0	"	"	"	"	"
2-Hexanone	U			50.0	"	"	"	"	"
Dibromochloromethane	U			20.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			20.0	"	"	"	"	"
Chlorobenzene	U			20.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			20.0	"	"	"	"	"
o-Xylene	U			20.0	"	"	"	"	"
Styrene	U			20.0	"	"	"	"	"
Bromoform	U			20.0	"	"	"	"	"
Isopropylbenzene	74.4			20.0	"	"	"	"	"
Bromobenzene	U			20.0	"	"	"	"	"
1,2,3-Trichloropropane	U			20.0	"	"	"	"	"





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77 West Jackson Boulevard  
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Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW003-200304 (2003006-05RE2)

Matrix: Water

Sampled: Mar-04-20 09:05

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
n-Propylbenzene	76.0			20.0	ug/L	10	B20D016	Mar-10-20	Mar-11-20
2-Chlorotoluene	U			20.0	"	"	"	"	"
1,3,5-Trimethylbenzene	121			20.0	"	"	"	"	"
4-Chlorotoluene	U			20.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			20.0	"	"	"	"	"
tert-Butylbenzene	U			20.0	"	"	"	"	"
1,2,4-Trimethylbenzene	329			20.0	"	"	"	"	"
sec-Butylbenzene	27.8			20.0	"	"	"	"	"
1,3-Dichlorobenzene	U			20.0	"	"	"	"	"
p-Isopropyltoluene	U			20.0	"	"	"	"	"
1,4-Dichlorobenzene	U			20.0	"	"	"	"	"
1,2-Dichlorobenzene	U			20.0	"	"	"	"	"
n-Butylbenzene	23.7	J		20.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			20.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			20.0	"	"	"	"	"
Hexachlorobutadiene	U			20.0	"	"	"	"	"
Naphthalene	29.1			20.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			20.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.48			94.2%		73-124	"	"	"
1,2-Dichloroethane-d4	10.6			105%		84-122	"	"	"
Toluene-d8	9.37			93.7%		88-108	"	"	"
4-Bromofluorobenzene	10.2			102%		84-108	"	"	"

A11-MW004A-200304 (2003006-06RE1)

Matrix: Water

Sampled: Mar-04-20 15:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Toluene	45300			2000	ug/L	1000	B20C019	Mar-10-20	Mar-10-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.61			95.5%		73-124	"	"	"
1,2-Dichloroethane-d4	10.9			108%		84-122	"	"	"
Toluene-d8	9.86			98.6%		88-108	"	"	"
4-Bromofluorobenzene	8.80			88.0%		84-108	"	"	"





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536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004A-200304 (2003006-06RE2)

Matrix: Water

Sampled: Mar-04-20 15:20

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			50.0	ug/L	25	B20D016	Mar-10-20	Mar-11-20
Chloromethane	U			50.0	"	"	"	"	"
Vinyl chloride	U			50.0	"	"	"	"	"
Bromomethane	U	(LCS), J		50.0	"	"	"	"	"
Chloroethane	U			50.0	"	"	"	"	"
Trichlorofluoromethane	U			50.0	"	"	"	"	"
1,1-Dichloroethene	U			50.0	"	"	"	"	"
Acetone	U			312	"	"	"	"	"
Carbon disulfide	U			50.0	"	"	"	"	"
Methylene chloride	U			50.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			50.0	"	"	"	"	"
1,1-Dichloroethane	U			50.0	"	"	"	"	"
2,2-Dichloropropane	U			50.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			50.0	"	"	"	"	"
2-Butanone	U			312	"	"	"	"	"
Bromochloromethane	U			50.0	"	"	"	"	"
Chloroform	U			50.0	"	"	"	"	"
1,1,1-Trichloroethane	U			50.0	"	"	"	"	"
Carbon tetrachloride	U			50.0	"	"	"	"	"
1,1-Dichloropropene	U			50.0	"	"	"	"	"
Benzene	U			50.0	"	"	"	"	"
1,2-Dichloroethane	U			50.0	"	"	"	"	"
Trichloroethene	U			50.0	"	"	"	"	"
1,2-Dichloropropane	U			50.0	"	"	"	"	"
Dibromomethane	U			50.0	"	"	"	"	"
Bromodichloromethane	U			50.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			50.0	"	"	"	"	"
4-Methyl-2-pentanone	U			125	"	"	"	"	"
trans-1,3-Dichloropropene	U			50.0	"	"	"	"	"
1,1,2-Trichloroethane	U			50.0	"	"	"	"	"
Tetrachloroethene	U			50.0	"	"	"	"	"
1,3-Dichloropropane	U			50.0	"	"	"	"	"
2-Hexanone	U			125	"	"	"	"	"
Dibromochloromethane	U			50.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			50.0	"	"	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

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Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004A-200304 (2003006-06RE2)

Matrix: Water

Sampled: Mar-04-20 15:20

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Chlorobenzene	U			50.0	ug/L	25	B20D016	Mar-10-20	Mar-11-20
1,1,1,2-Tetrachloroethane	U			50.0	"	"	"	"	"
Ethylbenzene	260			50.0	"	"	"	"	"
m+p-Xylene	414			100	"	"	"	"	"
o-Xylene	U			50.0	"	"	"	"	"
Styrene	U			50.0	"	"	"	"	"
Bromoform	U			50.0	"	"	"	"	"
Isopropylbenzene	U			50.0	"	"	"	"	"
Bromobenzene	U			50.0	"	"	"	"	"
1,2,3-Trichloropropane	U			50.0	"	"	"	"	"
n-Propylbenzene	U			50.0	"	"	"	"	"
2-Chlorotoluene	U			50.0	"	"	"	"	"
1,3,5-Trimethylbenzene	U			50.0	"	"	"	"	"
4-Chlorotoluene	U			50.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			50.0	"	"	"	"	"
tert-Butylbenzene	U			50.0	"	"	"	"	"
1,2,4-Trimethylbenzene	U			50.0	"	"	"	"	"
sec-Butylbenzene	U			50.0	"	"	"	"	"
1,3-Dichlorobenzene	U			50.0	"	"	"	"	"
p-Isopropyltoluene	U			50.0	"	"	"	"	"
1,4-Dichlorobenzene	U			50.0	"	"	"	"	"
1,2-Dichlorobenzene	U			50.0	"	"	"	"	"
n-Butylbenzene	U			50.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			50.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			50.0	"	"	"	"	"
Hexachlorobutadiene	U			50.0	"	"	"	"	"
Naphthalene	U			50.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			50.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.57			95.2%		73-124	"	"	"
1,2-Dichloroethane-d4	11.2			111%		84-122	"	"	"
Toluene-d8	9.86			98.6%		88-108	"	"	"
4-Bromofluorobenzene	9.19			91.9%		84-108	"	"	"





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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
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Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

### Volatiles by GC/MS, EPA 8260C (modified)

### US EPA Region 5 LSASD Analytical Services Branch

A11-MW004B-200303 (2003006-07RE2)

Matrix: Water

Sampled: Mar-03-20 16:05

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20D016	Mar-10-20	Mar-11-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U	(LCS), J		2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	5.86			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	5.29			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"





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77 West Jackson Boulevard  
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Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004B-200303 (2003006-07RE2)

Matrix: Water

Sampled: Mar-03-20 16:05

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2-Dibromoethane (EDB)	U			2.00	ug/L	1	B20D016	Mar-10-20	Mar-11-20
Chlorobenzene	U			2.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.3			102%		73-124	"	"	"
1,2-Dichloroethane-d4	11.4			113%		84-122	"	"	"
Toluene-d8	9.83			98.3%		88-108	"	"	"
4-Bromofluorobenzene	8.72			87.2%		84-108	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW005-200303 (2003006-08RE2)

Matrix: Water

Sampled: Mar-03-20 13:40

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20D016	Mar-10-20	Mar-11-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U	(LCS), J		2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U	(MS), L		2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	3.77			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	2.92			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"





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Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW005-200303 (2003006-08RE2)

Matrix: Water

Sampled: Mar-03-20 13:40

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2-Dibromoethane (EDB)	U			2.00	ug/L	1	B20D016	Mar-10-20	Mar-11-20
Chlorobenzene	U			2.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.2			101%		73-124	"	"	"
1,2-Dichloroethane-d4	11.5			114%		84-122	"	"	"
Toluene-d8	9.56			95.6%		88-108	"	"	"
4-Bromofluorobenzene	8.61			86.1%		84-108	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW006-200303 (2003006-09RE2)

Matrix: Water

Sampled: Mar-03-20 11:00

Received: Mar-05-20 14:45

Sample Qualifiers: (LCS), J

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20D016	Mar-10-20	Mar-11-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U	(LCS), J		2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	2.62			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"





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536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW006-200303 (2003006-09RE2)**

**Matrix: Water**

**Sampled: Mar-03-20 11:00**

**Received: Mar-05-20 14:45**

**Sample Qualifiers: (LCS), J**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2-Dibromoethane (EDB)	U			2.00	ug/L	1	B20D016	Mar-10-20	Mar-11-20
Chlorobenzene	U			2.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.85			98.0%		73-124	"	"	"
1,2-Dichloroethane-d4	11.1			110%		84-122	"	"	"
Toluene-d8	9.35			93.5%		88-108	"	"	"
4-Bromofluorobenzene	8.88			88.8%		84-108	"	"	"





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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200304 (2003006-10)

Matrix: Water

Sampled: Mar-04-20 11:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethylbenzene	959			100	ug/L	50	B20C012	Mar-06-20	Mar-09-20
m+p-Xylene	3050			200	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.9			109%		73-124	"	"	"
1,2-Dichloroethane-d4	11.0			110%		84-122	"	"	"
Toluene-d8	10.3			103%		88-108	"	"	"
4-Bromofluorobenzene	9.67			96.7%		84-108	"	"	"

A11-MW007-200304 (2003006-10RE1)

Matrix: Water

Sampled: Mar-04-20 11:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			4.00	ug/L	2	B20C019	Mar-10-20	Mar-11-20
Chloromethane	U			4.00	"	"	"	"	"
Vinyl chloride	U			4.00	"	"	"	"	"
Bromomethane	U			4.00	"	"	"	"	"
Chloroethane	U			4.00	"	"	"	"	"
Trichlorofluoromethane	U			4.00	"	"	"	"	"
1,1-Dichloroethene	U			4.00	"	"	"	"	"
Acetone	U			25.0	"	"	"	"	"
Carbon disulfide	U			4.00	"	"	"	"	"
Methylene chloride	U			4.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			4.00	"	"	"	"	"
1,1-Dichloroethane	U			4.00	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		4.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			4.00	"	"	"	"	"
2-Butanone	U			25.0	"	"	"	"	"
Bromochloromethane	U			4.00	"	"	"	"	"
Chloroform	U			4.00	"	"	"	"	"
1,1,1-Trichloroethane	U			4.00	"	"	"	"	"
Carbon tetrachloride	U			4.00	"	"	"	"	"
1,1-Dichloropropene	U			4.00	"	"	"	"	"
Benzene	U			4.00	"	"	"	"	"
1,2-Dichloroethane	U			4.00	"	"	"	"	"
Trichloroethene	U			4.00	"	"	"	"	"
1,2-Dichloropropane	U			4.00	"	"	"	"	"
Dibromomethane	U			4.00	"	"	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW007-200304 (2003006-10RE1)**

**Matrix: Water**

**Sampled: Mar-04-20 11:20**

**Received: Mar-05-20 14:45**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Bromodichloromethane	U			4.00	ug/L	2	B20C019	Mar-10-20	Mar-11-20
cis-1,3-Dichloropropene	U			4.00	"	"	"	"	"
4-Methyl-2-pentanone	U			10.0	"	"	"	"	"
Toluene	U			4.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			4.00	"	"	"	"	"
1,1,2-Trichloroethane	U			4.00	"	"	"	"	"
Tetrachloroethene	U			4.00	"	"	"	"	"
1,3-Dichloropropane	U			4.00	"	"	"	"	"
2-Hexanone	U			10.0	"	"	"	"	"
Dibromochloromethane	U			4.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			4.00	"	"	"	"	"
Chlorobenzene	U			4.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			4.00	"	"	"	"	"
o-Xylene	U			4.00	"	"	"	"	"
Styrene	U			4.00	"	"	"	"	"
Bromoform	U			4.00	"	"	"	"	"
Isopropylbenzene	12.0			4.00	"	"	"	"	"
Bromobenzene	U			4.00	"	"	"	"	"
1,2,3-Trichloropropane	U			4.00	"	"	"	"	"
n-Propylbenzene	6.68			4.00	"	"	"	"	"
2-Chlorotoluene	U			4.00	"	"	"	"	"
1,3,5-Trimethylbenzene	4.84			4.00	"	"	"	"	"
4-Chlorotoluene	U			4.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			4.00	"	"	"	"	"
tert-Butylbenzene	U			4.00	"	"	"	"	"
1,2,4-Trimethylbenzene	22.6			4.00	"	"	"	"	"
sec-Butylbenzene	U			4.00	"	"	"	"	"
1,3-Dichlorobenzene	U			4.00	"	"	"	"	"
p-Isopropyltoluene	U			4.00	"	"	"	"	"
1,4-Dichlorobenzene	U			4.00	"	"	"	"	"
1,2-Dichlorobenzene	U			4.00	"	"	"	"	"
n-Butylbenzene	U			4.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			4.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			4.00	"	"	"	"	"
Hexachlorobutadiene	U			4.00	"	"	"	"	"
Naphthalene	U			4.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			4.00	"	"	"	"	"





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536 South Clark Street, Chicago, IL 60605  
Phone: (312) 353-8370 Fax: (312) 886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW007-200304 (2003006-10RE1)**

**Matrix: Water**

**Sampled: Mar-04-20 11:20**

**Received: Mar-05-20 14:45**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>Dibromofluoromethane</i>	9.74			96.8%		73-124	B20C019	Mar-10-20	Mar-11-20
<i>1,2-Dichloroethane-d4</i>	10.7			106%		84-122	"	"	"
<i>Toluene-d8</i>	9.63			96.3%		88-108	"	"	"
<i>4-Bromofluorobenzene</i>	10.6			106%		84-108	"	"	"

**A11-MW007-200304-D (2003006-11)**

**Matrix: Water**

**Sampled: Mar-04-20 11:20**

**Received: Mar-05-20 14:45**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Ethylbenzene</b>	<b>863</b>			100	ug/L	50	B20C012	Mar-06-20	Mar-09-20
<b>m+p-Xylene</b>	<b>2800</b>			200	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>Dibromofluoromethane</i>	10.7			106%		73-124	"	"	"
<i>1,2-Dichloroethane-d4</i>	11.1			110%		84-122	"	"	"
<i>Toluene-d8</i>	10.4			104%		88-108	"	"	"
<i>4-Bromofluorobenzene</i>	9.80			98.0%		84-108	"	"	"

**A11-MW007-200304-D (2003006-11RE1)**

**Matrix: Water**

**Sampled: Mar-04-20 11:20**

**Received: Mar-05-20 14:45**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Dichlorodifluoromethane</b>	U			4.00	ug/L	2	B20C019	Mar-10-20	Mar-10-20
<b>Chloromethane</b>	U			4.00	"	"	"	"	"
<b>Vinyl chloride</b>	U			4.00	"	"	"	"	"
<b>Bromomethane</b>	U			4.00	"	"	"	"	"
<b>Chloroethane</b>	U			4.00	"	"	"	"	"
<b>Trichlorofluoromethane</b>	U			4.00	"	"	"	"	"
<b>1,1-Dichloroethene</b>	U			4.00	"	"	"	"	"
<b>Acetone</b>	U			25.0	"	"	"	"	"
<b>Carbon disulfide</b>	U			4.00	"	"	"	"	"
<b>Methylene chloride</b>	U			4.00	"	"	"	"	"
<b>trans-1,2-Dichloroethene</b>	U			4.00	"	"	"	"	"
<b>1,1-Dichloroethane</b>	U			4.00	"	"	"	"	"
<b>2,2-Dichloropropane</b>	U	(LCS), J		4.00	"	"	"	"	"
<b>cis-1,2-Dichloroethene</b>	U			4.00	"	"	"	"	"
<b>2-Butanone</b>	U			25.0	"	"	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200304-D (2003006-11RE1)

Matrix: Water

Sampled: Mar-04-20 11:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Bromochloromethane	U			4.00	ug/L	2	B20C019	Mar-10-20	Mar-10-20
Chloroform	U			4.00	"	"	"	"	"
1,1,1-Trichloroethane	U			4.00	"	"	"	"	"
Carbon tetrachloride	U			4.00	"	"	"	"	"
1,1-Dichloropropene	U			4.00	"	"	"	"	"
Benzene	U			4.00	"	"	"	"	"
1,2-Dichloroethane	U			4.00	"	"	"	"	"
Trichloroethene	U			4.00	"	"	"	"	"
1,2-Dichloropropane	U			4.00	"	"	"	"	"
Dibromomethane	U			4.00	"	"	"	"	"
Bromodichloromethane	U			4.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			4.00	"	"	"	"	"
4-Methyl-2-pentanone	U			10.0	"	"	"	"	"
Toluene	U			4.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			4.00	"	"	"	"	"
1,1,2-Trichloroethane	U			4.00	"	"	"	"	"
Tetrachloroethene	U			4.00	"	"	"	"	"
1,3-Dichloropropane	U			4.00	"	"	"	"	"
2-Hexanone	U			10.0	"	"	"	"	"
Dibromochloromethane	U			4.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			4.00	"	"	"	"	"
Chlorobenzene	U			4.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			4.00	"	"	"	"	"
o-Xylene	U			4.00	"	"	"	"	"
Styrene	U			4.00	"	"	"	"	"
Bromoform	U			4.00	"	"	"	"	"
Isopropylbenzene	11.0			4.00	"	"	"	"	"
Bromobenzene	U			4.00	"	"	"	"	"
1,2,3-Trichloropropane	U			4.00	"	"	"	"	"
n-Propylbenzene	6.03			4.00	"	"	"	"	"
2-Chlorotoluene	U			4.00	"	"	"	"	"
1,3,5-Trimethylbenzene	4.57			4.00	"	"	"	"	"
4-Chlorotoluene	U			4.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			4.00	"	"	"	"	"
tert-Butylbenzene	U			4.00	"	"	"	"	"
1,2,4-Trimethylbenzene	21.0			4.00	"	"	"	"	"
sec-Butylbenzene	U			4.00	"	"	"	"	"





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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Apr-20-20 17:34

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200304-D (2003006-11RE1)

Matrix: Water

Sampled: Mar-04-20 11:20

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,3-Dichlorobenzene	U			4.00	ug/L	2	B20C019	Mar-10-20	Mar-10-20
p-Isopropyltoluene	U			4.00	"	"	"	"	"
1,4-Dichlorobenzene	U			4.00	"	"	"	"	"
1,2-Dichlorobenzene	U			4.00	"	"	"	"	"
n-Butylbenzene	U			4.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			4.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			4.00	"	"	"	"	"
Hexachlorobutadiene	U			4.00	"	"	"	"	"
Naphthalene	U			4.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			4.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.2			101%		73-124	"	"	"
1,2-Dichloroethane-d4	10.9			109%		84-122	"	"	"
Toluene-d8	9.90			99.0%		88-108	"	"	"
4-Bromofluorobenzene	10.7			107%		84-108	"	"	"

A11-TB001-200303 (2003006-12)

Matrix: Water

Sampled: Mar-03-20 08:00

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	2.33			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-TB001-200303 (2003006-12)

Matrix: Water

Sampled: Mar-03-20 08:00

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1-Trichloroethane	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"

Report Name: 2003006 VOA - 8260 FINAL Apr 20 20 1734





## Environmental Protection Agency Region 5

### US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

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77 West Jackson Boulevard  
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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Apr-20-20 17:34

### Volatiles by GC/MS, EPA 8260C (modified)

### US EPA Region 5 LSASD Analytical Services Branch

A11-TB001-200303 (2003006-12)

Matrix: Water

Sampled: Mar-03-20 08:00

Received: Mar-05-20 14:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,3-Dichlorobenzene	U			2.00	ug/L	1	B20C012	Mar-06-20	Mar-09-20
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.7			106%		73-124	"	"	"
1,2-Dichloroethane-d4	10.5			104%		84-122	"	"	"
Toluene-d8	10.4			104%		88-108	"	"	"
4-Bromofluorobenzene	9.22			92.2%		84-108	"	"	"



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 20030087  
**Laboratory:** STAT Analysis Corporation / Eurofins Test America  
**Matrix:** Groundwater  
**Collection date:** 03/03/20  
**Analysis/Methods:**

Wet Chemistry:  
     Anions 300.0  
     Alkalinity M2320 B  
     Dissolved Gases - Methane - RSK-175

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>
20030087-01	A11-MW001-200303
20030087-02	A11-MW005-200303
20030087-03	A11-MW006-200303
20030087-04	A11-MW004B-200303

Data validation was performed in accordance with the specific analytical methods, the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017), National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?			N/A
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)		Yes	
Laboratory Control Spike Duplicates RPD within limits?		Yes	
Laboratory Duplicate RPDs within limits?		N/A	
<u>Comments (note deviations):</u>			

<u>Field Duplicates</u>	<u>Sample</u>	<u>Duplicate</u>	<u>%RPD</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					

<u>MS/MSD</u>	<u>%RPD</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate 20030087-001BMS/BMSD (20030087-01)	Acceptable	20%		
Sulfate 20030087-001BMS/BMSD (20030087-01)	Acceptable	20%		
Alkalinity 20030087-001BMS/BMSD	Acceptable	20%		

<u>LCS/LCSD</u>	<u>%RPD</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				

<u>Laboratory Duplicate</u>	<u>%RPD</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	No		
Were the Field Blanks results all < RL?	N/A		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			



<b>Blanks</b>		<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Nitrogen, Nitrate ICMBW1 030720	Nitrogen	0.048 J	0.2	None	Sample results > RL
Sulfate ICMBW1 030720	Sulfate	0.395 J	4.0	None	Sample results > RL
Alkalinity ALKMBW1 030720		Nondetect			
<b>ICB/CCB</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
ICB	Nitrogen, Nitrate	0.05	0.2	None	Sample results > RL
ICB	Sulfate	0.363	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.05	0.2	None	Sample results > RL
CCB	Sulfate	0.379	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.051	0.2	None	Sample results > RL
CCB	Sulfate	0.381	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.054	0.2	None	Sample results > RL
CCB	Sulfate	0.365	4.0	None	Sample results > RL
<b>Field Blank</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A					
<b>Surrogates</b>		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A					
<b>MS/MSD</b>		<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Nitrogen, Nitrate 20030087-001BMS/BMSD (20030087-01)		Acceptable	90-110		
Sulfate 20030087-001BMS/BMSD (20030087-01)		Acceptable	90-110		
Alkalinity 17030290-003BMS/MSD		Acceptable	75-125		
<b>LCS/LCSD</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Nitrogen, Nitrate ICLCSW1 030720		Acceptable	90-110		
Sulfate ICLCSW1 030720		Acceptable	90-110		
Alkalinity ALKLCSW1 030720		Acceptable	90-110		
<b>ICV</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
March 7 - 21:45	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			
<b>CCV</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
March 7 - 21:32	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			
March 8 - 00:23	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			



Tune  
N/A

Internal Standards  
N/A

Area

Area Lower / Upper  
Limit

Qualifiers Associated Samples



**Methane (RSK-175)****Precision:**

Are the field duplicate relative percent differences (RPD)  $\leq 30\%$  (aqueous)?  
 Were the Matrix Spike Duplicate RPDs  $\leq 20\%$ ? (Or lab defined limits)  
 Laboratory Control Spike Duplicates RPD within limits?  
 Laboratory Duplicate RPDs within limits?

Yes	No	N/A
		N/A
	Yes	
	Yes	
	N/A	

Comments (note deviations):

Field Duplicates	Sample	Duplicate	%RPD	Qualifiers	Associated Samples
N/A					

MS/MSD	%RPD	Limit	Qualifiers	Associated Samples
Methane				
608-181174-1 MS/MSD	Acceptable			

LCS/LCSD	%RPD	Limits	Qualifiers	Associated Samples
Methane				
LCS 680-610346 / 3 / 4	Acceptable			
LCS 680-611124 / 6 / 7	Acceptable			

Laboratory Duplicate	%RPD	Limits	Qualifiers	Associated Samples
N/A				

**Accuracy:**

Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency  $\geq 5\%$  and laboratory determined control limits)  
 Laboratory Control Sample criteria met?  
 Were the Laboratory Method Blank results all  $< RL$ ?  
 Were the Field Blanks results all  $< RL$ ?  
 Was the ICAL criteria met?  
 Was the CCV criteria met?  
 Was the Tuning criteria met?  
 Were the Surrogate % recoveries within laboratory determined control limits?  
 Were the Internal Standard areas within  $\pm 50 - 150\%$ ?

Yes	No	N/A
	Yes	
	Yes	
	Yes	
	N/A	
	Yes	
	Yes	
	N/A	
	N/A	
	N/A	

Comments (note deviations):

Blanks	Concentration (mg/L)	MDL / PQL	Qualifiers	Associated Samples
Methane				
MB 680-610346 / 8	Nondetect			
MB 680-611124 / 8	Nondetect			

Field Blank	Concentration	MDL / PQL	Qualifiers	Associated Samples
N/A				

Surrogates	%R	Limit	Qualifiers	Associated Samples
N/A				

MS/MSD	%R	Limits (%)	Qualifiers	Associated Samples
Methane				
608-181174-1 MS/MSD	Acceptable			

LCS/LCSD	%R	Limits	Qualifiers	Associated Samples
Methane				
LCS 680-610346 / 3 / 4	Acceptable			
LCS 680-611124 / 6 / 7	Acceptable			

ICAL	RRF	%RSD	Qualifiers	Associated Samples
3/4/2020 9:12	Acceptable	Acceptable		



<b>CCV</b>	<b>RRF</b>	<b>%D</b>	<b>Limits</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
3/10/2020 15:13	Acceptable	Acceptable			
3/10/2020 19:01	Acceptable	Acceptable			

**Tune**  
N/A

<b>Internal Standards</b>	<b>Area</b>	<b>Area Lower / Upper Limit</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
N/A				

**Representativeness:**

Were sampling procedures and design criteria met?

Were holding times met?

Was preservation criteria met? (0° C - 6° C)

Were Chain-of-Custody records complete and provided in data package?

Comments (note deviations): The cooler temperatures were 3.1 and 1.6 ° C.

**Yes No N/A**

Yes

Yes

Yes

Yes

<b>Preservation</b>	<b>Cooler Temperature (Degrees C)</b>	<b>Preservation Criteria</b>	<b>Qualifier</b>	<b>Associated Samples</b>
	Acceptable			

<b>Holding Times</b>	<b>Analyte</b>	<b>Days to Extraction</b>	<b>HT Criteria</b>	<b>Qualifier</b>	<b>Associated Samples</b>
		Acceptable			

**Comparability:**

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

**Yes No N/A**

Yes

**Completeness (90%):**

Are all data in this SDG usable?

Comments (note deviations):

**Yes No N/A**

Yes

**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):

**Yes No N/A**

Yes

Yes

**Comment:**

Data is usable as reported.

Data Validator:

*Kristine Molloy*

Date: 12/5/2020

Data Reviewer:

Cherie Zakowski

Date: 12/8/2021



**STAT Analysis Corporation**

2242 West Harrison St., Suite 200, Chicago, IL 60612-3766

Tel: (312) 733-0551 Fax: (312) 733-2386 STATinfo@STATAnalysis.com

Accreditations: IEPA ELAP 100445; ORELAP IL300001; AIHA-LAP, LLC 101160; NVLAP LabCode 101202-0

Date Reported: March 30, 2020

Date Printed: March 30, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Semi Annual GW Sampli Work Order: 20030087 Revision 0

Lab ID: 20030087-001

Collection Date: 3/3/2020 9:20:00 AM

Client Sample ID A11-MW001-200303

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>				Prep Date: 3/7/2020	Analyst: MD
Nitrogen, Nitrate (As N)	3.5	0.20	*	mg/L	1	3/7/2020
Sulfate	45	4.0	*	mg/L	1	3/7/2020
<b>Alkalinity</b>	<b>M2320 B</b>				Prep Date: 3/7/2020	Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	360	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>				Prep Date:	Analyst: SUB
Methane	ND	0.00058		mg/L	1	3/10/2020

Lab ID: 20030087-002

Collection Date: 3/3/2020 1:40:00 PM

Client Sample ID A11-MW005-200303

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>				Prep Date: 3/7/2020	Analyst: MD
Nitrogen, Nitrate (As N)	2.5	0.20	*	mg/L	1	3/7/2020
Sulfate	31	4.0	*	mg/L	1	3/7/2020
<b>Alkalinity</b>	<b>M2320 B</b>				Prep Date: 3/7/2020	Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	320	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>				Prep Date:	Analyst: SUB
Methane	ND	0.00058		mg/L	1	3/10/2020

Lab ID: 20030087-003

Collection Date: 3/3/2020 11:00:00 AM

Client Sample ID A11-MW006-200303

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>				Prep Date: 3/7/2020	Analyst: MD
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	3/7/2020
Sulfate	35	4.0	*	mg/L	1	3/7/2020
<b>Alkalinity</b>	<b>M2320 B</b>				Prep Date: 3/7/2020	Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	490	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>				Prep Date:	Analyst: SUB
Methane	3.3	0.39		mg/L	1	3/16/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

HT - Sample received past holding time

\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H - Holding time exceeded



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Date Reported: March 30, 2020

Date Printed: March 30, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Semi Annual GW Sampli Work Order: 20030087 Revision 0

Lab ID: 20030087-004

Collection Date: 3/3/2020 4:05:00 PM

Client Sample ID A11-MW004B-200303

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>				Prep Date: <b>3/7/2020</b>	Analyst: <b>MD</b>
Nitrogen, Nitrate (As N)	1.1	0.20	*	mg/L	1	3/7/2020
Sulfate	19	4.0	*	mg/L	1	3/7/2020
<b>Alkalinity</b>	<b>M2320 B</b>				Prep Date: <b>3/7/2020</b>	Analyst: <b>MD</b>
Alkalinity, Total (As CaCO3)	330	200		mg/L CaCO3	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>				Prep Date:	Analyst: <b>SUB</b>
Methane	ND	0.00058		mg/L	1	3/10/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
HT - Sample received past holding time  
\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H - Holding time exceeded



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 20030133  
**Laboratory:** STAT Analysis Corporation / Eurofins Test America  
**Matrix:** Groundwater  
**Collection date:** 03/04/20  
**Analysis/Methods:**

Wet Chemistry:  
     Anions 300.0  
     Alkalinity M2320 B  
     Dissolved Gases - Methane - RSK-175

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>
20030133-001	A11-MW003-200304
20030133-002	A11-MW002-200304
20030133-003	A11-MW007-200304
20030133-004	A11-MW004A-200304
20030133-005	A11-MW004A-200304-D

Data validation was performed in accordance with the specific analytical methods, the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017), and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	Yes		
Laboratory Control Spike Duplicates RPD within limits?	N/A		
Laboratory Duplicate RPDs within limits?	N/A		
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b><u>Sample</u> A11-MW004A- 200304</b>	<b><u>Duplicate</u> A11-MW004A- 200304-D</b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
			Acceptable		

<b>MS/MSD</b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Nitrogen, Nitrate 20030087-001BMS/BMSD	Acceptable	20%		
Sulfate 20030087-001BMS/BMSD	Acceptable	20%		
Alkalinity 20030087-001BMS/BMSD	Acceptable	20%		

<b>LCS/LCSD</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	No		
Were the Field Blanks results all < RL?	N/A		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			



<b>Blanks</b>		<u>Concentration</u>	<u>MDL / PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate ICMBW1 030720	Nitrogen	0.048 J	0.2	None	Sample results > RL
Sulfate ICMBW1 030720	Sulfate	0.395 J	4.0	None	Sample results > RL
Alkalinity ALKMBW1 030720		Nondetect			
<b>ICB/CCB</b>		<u>Concentration</u>	<u>MDL / PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
ICB	Nitrogen, Nitrate	0.05	0.2	None	Sample results > RL
ICB	Sulfate	0.363	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.05	0.2	None	Sample results > RL
CCB	Sulfate	0.379	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.051	0.2	None	Sample results > RL
CCB	Sulfate	0.381	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.05	0.2	None	Sample results > RL
CCB	Sulfate	0.359	4.0	None	Sample results > RL
<b>Field Blank</b>		<u>Concentration</u>	<u>MDL / PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<b>Surrogates</b>		<u>%R</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<b>MS/MSD</b>		<u>%R</u>	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate 20030087-001BMS/BMSD		Acceptable	90-110		
Sulfate 20030087-001BMS/BMSD		Acceptable	90-110		
Alkalinity 20030087-001BMS/BMSD		Acceptable	75-125		
<b>LCS/LCSD</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate ICLCSW1 030720		Acceptable	90-110		
Sulfate ICLCSW1 030720		Acceptable	90-110		
Alkalinity ALKLCSW1 030720		Acceptable	80-120		
<b>ICV</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
March 7 - 21:45	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			
<b>CCV</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
March 7 - 21:32	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			
March 8 - 00:23	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			
March 8 - 04:02	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			



Tune  
N/A

Internal Standards  
N/A

Area

Area Lower / Upper  
Limit

Qualifiers Associated Samples



**Methane (RSK-175)**

<b>Precision:</b>					<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?					<b>Yes</b>		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)					<b>N/A</b>		
Laboratory Control Spike Duplicates RPD within limits?					<b>Yes</b>		
Laboratory Duplicate RPDs within limits?					<b>N/A</b>		
<u>Comments (note deviations):</u>							

<b>Field Duplicates</b>	<b><u>Sample</u> A11-MW004A- 200304</b>	<b><u>Duplicate</u> A11-MW004A- 200304-D</b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
			Acceptable		

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<b>MS/MSD</b> N/A	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
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<b>LCS/LCSD</b> Methane LCS 680-611285/ 3 / 4 LCS 680-611285/ 6 / 7	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	Acceptable			
	Acceptable			

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<b>Laboratory Duplicate</b> N/A	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
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<b>Accuracy:</b>					<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)							<b>N/A</b>
Laboratory Control Sample criteria met?					<b>Yes</b>		
Were the Laboratory Method Blank results all < RL?					<b>Yes</b>		
Were the Field Blanks results all < RL?					<b>N/A</b>		
Was the ICAL criteria met?					<b>Yes</b>		
Was the CCV criteria met?					<b>Yes</b>		
Was the Tuning criteria met?					<b>N/A</b>		
Were the Surrogate % recoveries within laboratory determined control limits?					<b>N/A</b>		
Were the Internal Standard areas within ± 50 - 150%?					<b>N/A</b>		
<u>Comments (note deviations):</u>							

<b>Blanks</b> Methane MB 680-611285/ 8	<b><u>Concentration</u> (mg/L)</b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	Nondetect			

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<b>Field Blank</b> N/A	<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
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<b>Surrogates</b> N/A	<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
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<b>MS/MSD</b> N/A	<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
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<b>LCS/LCSD</b> Methane LCS 680-611285/ 3 / 4 LCS 680-611285/ 6 / 7	<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	Acceptable			
	Acceptable			

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<b>ICAL</b> 2/17/2020 8:45 3/04/2020 9:12	<b><u>RRF</u></b>	<b><u>%RSD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	Acceptable	Acceptable		
	Acceptable	Acceptable		



CCV	RRF	%D	Limits	Qualifiers	Associated Samples
3/17/2020 17:11	Acceptable	Acceptable			
3/17/2020 17:37	Acceptable	Acceptable			
3/17/2020 20:35	Acceptable	Acceptable			
3/17/2020 8:38	Acceptable	Acceptable			

Tune  
N/A

Internal Standards	Area	Area Lower / Upper Limit	Qualifiers	Associated Samples
N/A				

#### Representativeness:

Were sampling procedures and design criteria met?

Were holding times met?

Was preservation criteria met? (0° C - 6° C)

Were Chain-of-Custody records complete and provided in data package?

Comments (note deviations): The cooler temperatures were 1.8 and 1.4 ° C.

Yes No N/A

Yes

Yes

Yes

Yes

Preservation	Cooler Temperature (Degrees C)	Preservation Criteria	Qualifier	Associated Samples
	Acceptable			

Holding Times	Analyte	Days to Extraction	HT Criteria	Qualifier	Associated Samples
		Acceptable			

#### Comparability:

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

Yes No N/A

Yes

#### Completeness (90%):

Are all data in this SDG usable?

Comments (note deviations):

Yes No N/A

Yes

#### Sensitivity:

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):

Yes No N/A

Yes

Yes

#### Comment:

Data is usable as reported.

Data Validator:

*Kristine Molloy*

Date: 1/22/2021

Data Reviewer:

Cherie Zakowski

Date: 1/25/2021



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2242 West Harrison St., Suite 200, Chicago, IL 60612-3766

Tel: (312) 733-0551 Fax: (312) 733-2386 STATinfo@STATAnalysis.com

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Date Reported: March 30, 2020

Date Printed: March 30, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Semi Annual GW Sampli Work Order: 20030133 Revision 0

Lab ID: 20030133-001

Collection Date: 3/4/2020 9:05:00 AM

Client Sample ID A11-MW003-200304

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 3/7/2020 Analyst: MD
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	3/7/2020
Sulfate	ND	4.0	*	mg/L	1	3/7/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 3/7/2020 Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	410	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	12	0.39		mg/L	1	3/17/2020

Lab ID: 20030133-002

Collection Date: 3/4/2020 1:35:00 PM

Client Sample ID A11-MW002-200304

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 3/7/2020 Analyst: MD
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	3/8/2020
Sulfate	ND	4.0	*	mg/L	1	3/8/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 3/7/2020 Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	370	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	16	0.39		mg/L	1	3/17/2020

Lab ID: 20030133-003

Collection Date: 3/4/2020 11:20:00 AM

Client Sample ID A11-MW007-200304

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 3/7/2020 Analyst: MD
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	3/8/2020
Sulfate	24	4.0	*	mg/L	1	3/8/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 3/7/2020 Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	330	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	5.5	0.39		mg/L	1	3/17/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

HT - Sample received past holding time

\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H - Holding time exceeded



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Date Reported: March 30, 2020

Date Printed: March 30, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Semi Annual GW Sampli Work Order: 20030133 Revision 0

Lab ID: 20030133-004

Collection Date: 3/4/2020 3:20:00 PM

Client Sample ID A11-MW004A-200304

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>				Prep Date: <b>3/7/2020</b>	Analyst: <b>MD</b>
Nitrogen, Nitrate (As N)	0.28	0.20	*	mg/L	1	3/8/2020
Sulfate	36	4.0	*	mg/L	1	3/8/2020
<b>Alkalinity</b>	<b>M2320 B</b>				Prep Date: <b>3/7/2020</b>	Analyst: <b>MD</b>
Alkalinity, Total (As CaCO <sub>3</sub> )	330	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>				Prep Date:	Analyst: <b>SUB</b>
Methane	0.51	0.00058		mg/L	1	3/17/2020

Lab ID: 20030133-005

Collection Date: 3/4/2020 3:20:00 PM

Client Sample ID A11-MW004A-200304-D

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>				Prep Date: <b>3/7/2020</b>	Analyst: <b>MD</b>
Nitrogen, Nitrate (As N)	0.38	0.20	*	mg/L	1	3/8/2020
Sulfate	36	4.0	*	mg/L	1	3/8/2020
<b>Alkalinity</b>	<b>M2320 B</b>				Prep Date: <b>3/7/2020</b>	Analyst: <b>MD</b>
Alkalinity, Total (As CaCO <sub>3</sub> )	320	200		mg/L CaCO <sub>3</sub>	1	3/7/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>				Prep Date:	Analyst: <b>SUB</b>
Methane	0.47	0.00058		mg/L	1	3/17/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
HT - Sample received past holding time  
\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H - Holding time exceeded



## **June 2020 Data Validation Reports and Data Packages**



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** E200604  
**Laboratory:** ESAT / TechLaw  
**Matrix:** Water  
**Collection date:** 6/9/2020 & 6/10/2020  
**Analysis/Methods:** 1,4-Dioxane - EPA 522 SIM

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
E200604-01	A11-TB001-200609	E200604-07	A11-MW003-200610
E200604-02	A11-MW006-200609	E200604-08	A11-MW007-200610
E200604-03	A11-MW005-200609	E200604-09	A11-MW007-200610-D
E200604-04	A11-MW001-200609	E200604-10	A11-MW002-200610
E200604-05	A11-MW004B-200609	E200604-11	A11-MW004A-200610
E200604-06	A11-FB001-200609		

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**1,4-Dioxane EPA 520 SIM**

Precision:	Yes	No	N/A
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	Yes		
Laboratory Control Spike Duplicates RPD within limits?	Yes		
Laboratory Duplicate RPDs within limits?	N/A		
Comments (note deviations):			

Field Duplicates	Sample A11-MW007-200610 ND	Duplicate A11-MW007-200610-D ND	%RPD Acceptable	Qualifiers	Associated Samples
MS/MSD E20F013-MS1 / MSD1	%RPD Acceptable	Limit		Qualifiers	Associated Samples
LCS/LCSD E20F013-BS1 / BSD1	%RPD Acceptable	Limits		Qualifiers	Associated Samples
Laboratory Duplicate N/A	%RPD	Limits		Qualifiers	Associated Samples

Accuracy:	Yes	No	N/A
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
MRL recoveries within criteria?	Yes		
Were the Field Blanks results all < RL?	Yes		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	Yes		
Were the Surrogate % recoveries within laboratory determined control limits?	Yes		
Were the Internal Standard areas within ± 50 - 150%?	Yes		
Comments (note deviations):			

Blanks	Concentration (µg/L)	MDL /RL	Qualifiers	Associated Samples
E20F013-BLK1	Nondetect			
Field Blank	Concentration	MDL /RL	Qualifiers	Associated Samples
A11-TB001-200609	Nondetect			
A11-FB001-200609	Nondetect			



Surrogates	<u>%R</u> Acceptable	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>MS/MSD</b> E20F013-MS1 / MSD1	<u>%R</u> Acceptable	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>LCS/LCSD</b> E20F013-BS1 / BSD1	<u>%R</u> Acceptable	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>MRL Check</b> E20F013-MRL1	<u>%R</u> Acceptable	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>ICAL</b> 5/ 27 / 2020 - 13:20	<u>RRF</u> Acceptable	<u>%RSD</u> Acceptable	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>ICV / CCV</b> ICV 5/ 27 / 2020 - 13:00	<u>RRF</u> Acceptable	<u>%D</u> Acceptable	<u>Qualifiers</u>	<u>Associated Samples</u>
CCV 6/ 24 / 2020 - 10:03 6/ 24 / 2020 - 5:42	Acceptable Acceptable	Acceptable Acceptable		
<b>Tune</b> Acceptable				

Internal Standards	<u>Area</u> Acceptable	<u>Area Lower / Upper Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
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**Representativeness:**

Were sampling procedures and design criteria met?

Were holding times met?

Was preservation criteria met? (0° C - 6° C)

Were Chain-of-Custody records complete and provided in data package?

Comments (note deviations): The cooler temperature was 4.6 ° C.

Yes No N/A

Yes

Yes

Yes

Yes

Preservation	<u>Cooler Temperature (Degrees C)</u> Acceptable	<u>Preservation Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
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Holding Times	<u>Analyte</u>	<u>Days to Analysis</u> Acceptable	<u>HT Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
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**Comparability:**

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):
Yes No N/A

Yes

**Completeness (90%):**

Are all data in this SDG usable?

Comments (note deviations):
Yes No N/A

Yes

**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):
Yes No N/A

Yes

Yes

**Comment:** Data is usable as reported.

Data Validator:

*Kristine Molloy*

Date: 1/22/2021

Data Reviewer:

*Cherie Zakowski*

Date: 1/25/2021



Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Howard Pham

Reported:  
Jul-16-20 13:56

## 1,4-Dioxane by GC-MS

### TechLaw - ESAT Contract

#### A11-TB001-200609 (E200604-01)

Matrix: Water

Sampled: Jun-09-20 08:00

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.207	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.918			88.8%		64-109	"	"	"

#### A11-MW006-200609 (E200604-02)

Matrix: Water

Sampled: Jun-09-20 11:00

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>7.53</b>			0.203	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	1.02			101%		64-109	"	"	"

#### A11-MW005-200609 (E200604-03)

Matrix: Water

Sampled: Jun-09-20 16:15

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>8.83</b>			0.205	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.977			95.4%		64-109	"	"	"

#### A11-MW001-200609 (E200604-04)

Matrix: Water

Sampled: Jun-09-20 08:40

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>14.1</b>			0.207	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	1.09			106%		64-109	"	"	"

#### A11-MW004B-200609 (E200604-05)

Matrix: Water

Sampled: Jun-09-20 13:05

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>11.7</b>			0.208	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	1.10			106%		64-109	"	"	"



Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Howard Pham

Reported:  
Jul-16-20 13:56

## 1,4-Dioxane by GC-MS

### TechLaw - ESAT Contract

#### A11-FB001-200609 (E200604-06)

Matrix: Water

Sampled: Jun-09-20 17:30

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.207	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	1.04			101%		64-109	"	"	"

#### A11-MW003-200610 (E200604-07)

Matrix: Water

Sampled: Jun-10-20 08:05

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>9.58</b>			0.205	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	1.03			100%		64-109	"	"	"

#### A11-MW007-200610 (E200604-08)

Matrix: Water

Sampled: Jun-10-20 09:50

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.205	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.943			92.0%		64-109	"	"	"

#### A11-MW007-200610-D (E200604-09)

Matrix: Water

Sampled: Jun-10-20 09:50

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.205	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.984			96.1%		64-109	"	"	"

#### A11-MW002-200610 (E200604-10)

Matrix: Water

Sampled: Jun-10-20 12:10

Received: Jun-11-20 10:18

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>4.03</b>			0.207	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	1.01			97.4%		64-109	"	"	"





TechLaw Inc ESAT Region 5  
536 South Clark Street, Suite 734  
Chicago, IL 60605  
(312) 353-8303  
(312) 353-5814 (Fax)  
[www.techlawinc.com](http://www.techlawinc.com)

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Howard Pham

**Reported:**  
Jul-16-20 13:56

**1,4-Dioxane by GC-MS**  
**TechLaw - ESAT Contract**

**A11-MW004A-200610 (E200604-11)**

**Matrix: Water**

**Sampled: Jun-10-20 14:25**

**Received: Jun-11-20 10:18**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>1.51</b>			0.203	ug/L	1	E20F013	Jun-23-20	Jun-24-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.915			90.0%		64-109	"	"	"



**Techlaw Document Controlled Number: 83074-8-33-704-DV-1330**  
UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V  
SUPERFUND AND EMERGENCY MANAGEMENT DIVISION

DATE:

SUBJECT: Review of Data  
Received for Review on: July 13, 2020

FROM: Allison Harvey, TechLaw Inc.  
Contractor, Environmental Services Assistance Team (ESAT)

THROUGH: Michelle Kerr  
Region 5 ESAT Contracting Officer's Representative

TO: Data User: CDM Smith  
Email Address: grabsjc@cdm.com

**This package was requested and reviewed as a Stage 4 Validation Electronic and Manual Deliverable (S4VEM)**

We have reviewed the data for the following case:

SITE Name: Southeast Rockford Groundwater Contamination (IL)

Case No: 48947 MA No: N/A SDG No: E3YF9

Number and Type of Samples: 11 waters (6 Trace Volatiles/ 5 L/M Volatiles)

Sample Numbers: E3YF9, E3YG0 – E3YG9

Laboratory: Chemtech Consulting Group (CHM) Hrs. for Review:

Following are our findings:



**Below is a summary of the out-of-control audits and the possible effects on the data for this case:**

Eleven (11) water samples were shipped to Chemtech Consulting Group (CHM) located in Mountainside, NJ. All samples were collected 06/09-10/2020 and received on 06/11/2020 intact and properly cooled. Six (6) samples; E3YF9 and E3YG0 thru E3YG4, were analyzed for the trace volatile analytes. Five (5) samples; E3YG5 thru E3YG9, were analyzed for the low level volatile analytes. All samples were analyzed according to CLP SOW SOM02.4, [Oct 2016] and reviewed according to the June 2010 Rev 1, March, 2014 Rev 2 QAPPs for Southeast Rockford Groundwater Contamination Site, the Illinois State QAPP, the September 2017 NFG for SOM02.4 (EPA-540-R-2017-002) and the Region 5 Organic CLP Validation SOP (DCN 83074-8-33-601-SO-1143.R1).

Samples E3YG1 and E3YG8 were designated by the samplers to be used for laboratory QC, i.e. MS/MSD analyses.

Sample E3YF9 was identified as a trip blank. Sample E3YG4 was identified as a field blank. Sample E3YG7 was identified as a field duplicate of sample E3YG6.

Only the qualifications reflected in the EXES Sample Summary report are described in this narrative.



**1. PRESERVATION AND HOLDING TIMES**

NONE FOUND.

**2. GC/MS and GC/ECD INSTRUMENT PERFORMANCE CHECK**

NONE FOUND.

**3. INITIAL CALIBRATION**

NONE FOUND.

**4. INITIAL CALIBRATION VERIFICATION**

NONE FOUND.

**5. CONTINUING CALIBRATION**

**Method – Volatile Organics**

EXES-1209

The following samples are associated with an opening or closing CCV with % Difference exceeding criteria. Detects are qualified as estimated J. Non-detects are qualified as estimated UJ.

E3YG5, E3YG6, E3YG7, E3YG8, E3YG8MS, E3YG8MSD, VBLK85  
Toluene

**6. BLANKS**

**Method – Trace Volatiles**

The following samples have analyte results reported less than CRQLs. The associated method blank results are less than CRQL. Detects are qualified U. Sample results have been reported at CRQLs.

VHBLK01  
Methylene chloride

The following samples have analyte results reported less than CRQLs. The associated trip blank (E3YF9) is less than CRQLs. Detects are qualified U. Sample results have been reported at CRQLs.

E3YG0, E3YG1, E3YG1MS, E3YG2, E3YG4  
Acetone



E3YG0, E3YG4  
cis-1,2-Dichloroethene

## **7. DEUTERATED MONITORING COMPOUNDS / SURROGATES**

### **Method – Trace Volatiles**

EXES-982

The following samples have DMC/surrogate percent recoveries greater than the primary maximum criteria. Detects are qualified as estimated J+. Non-detects are not qualified.

E3YF9  
Acetone, 2-Butanone, 4-Methyl-2-pentanone, 2-Hexanone

E3YG1, E3YG1MS, E3YG2, E3YG3, E3YG4  
4-Methyl-2-pentanone, 2-Hexanone

E3YG1MSD  
4-Methyl-2-pentanone, 2-Hexanone, Chlorobenzene, 1,3-Dichlorobenzene,  
1,4-Dichlorobenzene, 1,2-Dichlorobenzene, 1,2,4-Trichlorobenzene,  
1,2,3-Trichlorobenzene

## **8. MATRIX SPIKE/MATRIX SPIKE DUPLICATE**

### **Method – Volatile Organics**

EXES-1217

The following matrix spike/matrix spike duplicate samples have percent recoveries less than the expanded minimum criteria. Detects in the unspiked sample are qualified as estimated J. Non-detects in the unspiked sample are qualified as unusable R.

E3YG8MS, E3YG8MSD  
Toluene

EXES-559

The relative percent difference (RPD) between the following matrix spike and matrix spike duplicate recoveries is outside criteria. Detects in the unspiked sample are qualified as estimated J. Non-detects in the unspiked sample are not qualified.

E3YG8MS, E3YG8MSD  
Toluene



## **9. CLEANUP PROCEDURES**

Not required for these analyses.

## **10. LABORATORY CONTROL SAMPLE**

Not required for these analyses.

## **11. INTERNAL STANDARD**

NONE FOUND.

## **12. TARGET ANALYTE QUANTITATION LIMIT**

### **Method – Trace Volatiles**

EXES-790

The following samples have analyte results greater than or equal to method detection limit (MDL) and below contract required quantitation limit (CRQL). Detects are qualified as estimated J.

E3YF9

Acetone, cis-1,2-Dichloroethene

E3YG0

Chloroethane, 1,1-Dichloroethane, Cyclohexane, Trichloroethene, Isopropylbenzene

E3YG1, E3YG1MS, E3GY1MSD

trans-1,2-Dichloroethene, Chloroform, Bromodichloromethane, Tetrachloroethene, Dibromochloromethane

E3YG2

trans-1,2-Dichloroethene

E3YG3

trans-1,2-Dichloroethene, Tetrachloroethene

E3YG4

Carbon disulfide, 2-Butanone

VBLK09

Methylene chloride



## Method – Volatile Organics

### EXES-790

The following samples have analyte results greater than or equal to method detection limit (MDL) and below contract required quantitation limit (CRQL). Detects are qualified as estimated J.

#### E3YG5

Chloroethane, 1,1,1-Trichloroethane, 4-Methyl-2-pentanone

#### E3YG6, E3YG7

Methylcyclohexane, Tetrachloroethene

#### E3YG8

Vinyl chloride, Acetone, 1,1,1-Trichloroethane, Trichloroethene

#### E3YG8MS, E3YG8MSD

Vinyl chloride, Acetone, 1,1,1-Trichloroethane, 1,2-Dichlorobenzene

#### E3YG9

1,1,1-Trichloroethane, Cyclohexane, Trichloroethene, Isopropylbenzene

## 13. TENTATIVELY IDENTIFIED COMPOUNDS

Not Validated

## 14. SYSTEM PERFORMANCE

No problems found.

## 15. FIELD QC SAMPLES

Sample E3YF9 was identified as a trip blank. Sample E3YG4 was identified as a field blank. Sample E3YG7 was identified as a field duplicate of sample E3YG6.

Results are summarized in the following table:

Sample Type:	Trip Blank	Field Blank
Sample #:	A11-TB001-200609	A11-FB001-200609
CLP Sample:	<b>E3YF9</b>	<b>E3YG4</b>
Location:	A11-TB001	A11-FB001
Collection Date/Time:	6/9/2020 08:00	6/9/2020 17:30
Units:	µg/L	µg/L
Acetone	1.7 J	



Carbon disulfide	ND	0.090 J
cis-1,2-Dichloroethene	0.29 J	
2-Butanone	ND	1.2 J
Associated field samples:	E3YG0, E3YG1 E3YG2, E3YG3, E3YG4	E3YG0, E3YG1, E3YG2, E3YG3

ND = Not Detected.

Sample Type:	Field Sample	Field Duplicate	
Sample #:	A11-MW007-200610	A11-MW007-200610-D	
CLP Sample:	<b>E3YG6</b>	<b>E3YG7</b>	
Location:	A11-MW007	A11-MW007	
Collection Date/Time:	6/10/2020 9:50	6/10/2020 9:50	RPDs
Units:	µg/L	µg/L	%
Dilution factor:	1.0	1.0	
Methylcyclohexane	2.7 J	2.6 J	3.8
Tetrachloroethene	1.0 J	0.89 J	12
Isopropylbenzene	6.5	6.5	
CLP Sample:	<b>E3YG6DL</b>	<b>E3YG7DL</b>	
Dilution factor:	100.0	100.0	
Ethylbenzene	820	810	1.2
m,p-Xylene	2600	2600	0.0

‘\*’ – RPD value  $\geq$  20%.

The detection of analytes with RPDs greater than 20% in the field duplicates are qualified as estimated J. Nondetects are qualified as estimated UJ.

## 16. SAMPLE RESULTS

The following trace volatile samples have analyte results greater than the upper limit of calibration range. The samples were not re-analyzed at dilution because they are QC samples. Detects are qualified as estimated J.

E3YG8MS, E3YG8MSD

Methylcyclohexane, Toluene, Ethylbenzene, o-Xylene, m,p-Xylene

## 17. QAPP COMPLIANCE

The analytical package fulfilled the QAPP QC components requirements identified in the Southeast Rockford Groundwater Contamination QAPP.



Validation Data Qualifier Sheet

<u>Qualifiers</u>	<u>Data Qualifier Definitions</u>
U	The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
J	The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
J+	The result is an estimated quantity, but the results may be biased high.
J-	The result is an estimated quantity, but the results may be biased low.
NJ	The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
UJ	The analyte was analyzed for, but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
R	The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.
C	The target Pesticide or Aroclor analyte identification has been confirmed by Gas Chromatograph/Mass Spectrometer (GC/MS).
X	The target Pesticide or Aroclor analyte identification was not confirmed when GC/MS analysis was performed.



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YF9	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location: A11-TB001	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 08:00:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	1.7	J+	ug/L	1.7	J	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	0.29	J	ug/L	0.29	J	1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG0	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location: A11-MW006	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 11:00:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.44	J	ug/L	0.44	J	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5	U	ug/L	1.3	J	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	0.11	J	ug/L	0.11	J	1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	0.5	U	ug/L	0.28	J	1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Cyclohexane	Target	0.35	J	ug/L	0.35	J	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Target	2.0		ug/L	2.0		1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Target	0.14	J	ug/L	0.14	J	1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.14	J	ug/L	0.14	J	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Indane	TIC	0.57	JN	ug/L	0.57	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
n-Butyl ether	TIC	12	JN	ug/L	12	JN	1.0	YES	NV
Total Alkanes	TIC	1.5	BN	ug/L	1.5	BN	1.0	YES	NV
Di-sec-butyl ether	TIC	0.72	JN	ug/L	0.72	JN	1.0	YES	NV
unknown-01	TIC	2.4	J	ug/L	2.4	J	1.0	YES	NV
Pentalene, octahydro-	TIC	0.58	JN	ug/L	0.58	JN	1.0	YES	NV
Ethane, 1-chloro-1,1-difluoro-	TIC	1.7	JN	ug/L	1.7	JN	1.0	YES	NV
4-Octanone, 5-hydroxy-3,6-dimethyl	TIC	1.3	JN	ug/L	1.3	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG1	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location: A11-MW005	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 16:15:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	1.1		ug/L	1.1		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5	U	ug/L	0.87	J	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.15	J	ug/L	0.15	J	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.4		ug/L	6.4		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	1.3		ug/L	1.3		1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.45	J	ug/L	0.45	J	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	4.5		ug/L	4.5		1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Target	0.89		ug/L	0.89		1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.40	J	ug/L	0.40	J	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.39	J	ug/L	0.39	J	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.18	J	ug/L	0.18	J	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG1MS	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location:	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 16:15:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Spike	5.6		ug/L	5.6		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5	U	ug/L	1.1	J	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.16	J	ug/L	0.16	J	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.4		ug/L	6.4		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	1.3		ug/L	1.3		1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.45	J	ug/L	0.45	J	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	4.4		ug/L	4.4		1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Spike	4.8		ug/L	4.8		1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Spike	5.6		ug/L	5.6		1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.39	J	ug/L	0.39	J	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Spike	4.7		ug/L	4.7		1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.36	J	ug/L	0.36	J	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.19	J	ug/L	0.19	J	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Spike	4.8		ug/L	4.8		1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG1MSD	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location:	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 16:15:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Spike	5.8		ug/L	5.8		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.16	J	ug/L	0.16	J	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.6		ug/L	6.6		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	1.4		ug/L	1.4		1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.48	J	ug/L	0.48	J	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	4.6		ug/L	4.6		1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Spike	5.0		ug/L	5.0		1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Spike	5.7		ug/L	5.7		1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.38	J	ug/L	0.38	J	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Spike	4.8		ug/L	4.8		1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.39	J	ug/L	0.39	J	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.21	J	ug/L	0.21	J	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Spike	5.0	J+	ug/L	5.0		1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG2	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location: A11-MW001	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 08:40:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	1.4		ug/L	1.4		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5	U	ug/L	1.1	J	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.17	J	ug/L	0.17	J	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	7.5		ug/L	7.5		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	1.4		ug/L	1.4		1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	8.9		ug/L	8.9		1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Target	2.5		ug/L	2.5		1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	1.0		ug/L	1.0		1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethane, 1-chloro-1,1-difluoro-	TIC	0.77	JN	ug/L	0.77	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG3	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location: A11-MW004B	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 13:05:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	0.95		ug/L	0.95		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Carbon disulfide	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.16	J	ug/L	0.16	J	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.3		ug/L	6.3		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	1.3		ug/L	1.3		1.0	YES	S4VEM
2-Butanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	5.5		ug/L	5.5		1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Target	1.4		ug/L	1.4		1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Target	1.6		ug/L	1.6		1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.40	J	ug/L	0.40	J	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG4	Method: Trace Volatiles	Matrix: Water	MA Number:
Sample Location: A11-FB001	pH: 1.0	Sample Date: 06/09/2020	Sample Time: 17:30:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Vinyl chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromomethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Acetone	Target	5	U	ug/L	3.0	J	1.0	YES	S4VEM
Carbon disulfide	Target	0.090	J	ug/L	0.090	J	1.0	YES	S4VEM
Methyl Acetate	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylene chloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	0.5	U	ug/L	0.19	J	1.0	YES	S4VEM
2-Butanone	Target	1.2	J	ug/L	1.2	J	1.0	YES	S4VEM
Bromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chloroform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Cyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Benzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Trichloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Methylcyclohexane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromodichloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Toluene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
2-Hexanone	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Dibromochloromethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Chlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Ethylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
o-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
m,p-Xylene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Styrene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Bromoform	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Isopropylbenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	0.50	U	ug/L	0.50	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG5	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location: A11-MW003	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 08:05:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	1.9	J	ug/L	1.9	J	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.9		ug/L	6.9		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	1.3	J	ug/L	1.3	J	1.0	YES	S4VEM
Cyclohexane	Target	7.8		ug/L	7.8		1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylcyclohexane	Target	59		ug/L	59		1.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	2.2	J	ug/L	2.2	J	1.0	YES	S4VEM
Toluene	Target	7.6	J	ug/L	7.6		1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Ethylbenzene	Target	430	J	ug/L	430	JD	100.0	YES	S4VEM
o-xylene	Target	5.2		ug/L	5.2		1.0	YES	S4VEM
m,p-Xylene	Target	5100		ug/L	5100	D	100.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	45		ug/L	45		1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene, 4-ethyl-1,2-dimethyl-	TIC	13	JN	ug/L	13	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Benzene, 1,2,4,5-tetramethyl-	TIC	9.1	JN	ug/L	9.1	JN	1.0	YES	NV
Benzene, propyl-	TIC	53	JN	ug/L	53	JN	1.0	YES	NV
Benzene, 1,3-diethyl-	TIC	6.0	JN	ug/L	6.0	JN	1.0	YES	NV
p-Cymene	TIC	18	JN	ug/L	18	JN	1.0	YES	NV
Total Alkanes	TIC	120	BN	ug/L	120	BN	1.0	YES	NV
Benzene, 2-ethyl-1,4-dimethyl-	TIC	20	JN	ug/L	20	JN	1.0	YES	NV
Benzaldehyde, 2-methyl-	TIC	2.6	JN	ug/L	2.6	JN	1.0	YES	NV
Naphthalene, 1,2,3,4-tetrahydro-	TIC	18	JN	ug/L	18	JN	1.0	YES	NV
Benzene, (2-methyl-1-propenyl)-	TIC	27	JN	ug/L	27	JN	1.0	YES	NV
Benzene, 1,2,3,4-tetramethyl-	TIC	20	JN	ug/L	20	JN	1.0	YES	NV
Benzene, (1-methyl-1-butenyl)-	TIC	2.8	JN	ug/L	2.8	JN	1.0	YES	NV
Benzene, 1-methyl-4-propyl-	TIC	14	JN	ug/L	14	JN	1.0	YES	NV
Cyclohexene, 3-methyl-	TIC	3.1	JN	ug/L	3.1	JN	1.0	YES	NV
Benzene, (2-methylpropyl)-	TIC	4.5	JN	ug/L	4.5	JN	1.0	YES	NV
Benzene, 1-methyl-3-(1-methylethyl)	TIC	4.2	JN	ug/L	4.2	JN	1.0	YES	NV
Benzene, 1,2,3-trimethyl-	TIC	85	JN	ug/L	85	JN	1.0	YES	NV
Benzene, 1,2,4-trimethyl-	TIC	220	JN	ug/L	220	JN	1.0	YES	NV
Benzeneacetaldehyde, .alpha.-methy	TIC	15	JN	ug/L	15	JN	1.0	YES	NV
Benzene, 1-ethyl-2-methyl-	TIC	130	JN	ug/L	130	JN	1.0	YES	NV
Benzene, 1-ethyl-3-methyl-	TIC	54	JN	ug/L	54	JN	1.0	YES	NV
Benzene, 1,2-diethyl-	TIC	23	JN	ug/L	23	JN	1.0	YES	NV
Pentalene, octahydro-	TIC	6.5	JN	ug/L	6.5	JN	1.0	YES	NV
Indan, 1-methyl-	TIC	2.5	JN	ug/L	2.5	JN	1.0	YES	NV
o-Cymene	TIC	25	JN	ug/L	25	JN	1.0	YES	NV
1-Hexadecyne	TIC	2.8	JN	ug/L	2.8	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG6	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location: A11-MW007	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 09:50:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Cyclohexane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylcyclohexane	Target	2.7	J	ug/L	2.7	J	1.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Toluene	Target	5.0	UJ	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	1.0	J	ug/L	1.0	J	1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Ethylbenzene	Target	820		ug/L	820	D	100.0	YES	S4VEM
o-xylene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
m,p-Xylene	Target	2600		ug/L	2600	D	100.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	6.5		ug/L	6.5		1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Mesitylene	TIC	24	JN	ug/L	24	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV
Benzene, 1,2,3-trimethyl-	TIC	11	JN	ug/L	11	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG7	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location: A11-MW007	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 09:50:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Cyclohexane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylcyclohexane	Target	2.6	J	ug/L	2.6	J	1.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Toluene	Target	5.0	UJ	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	0.89	J	ug/L	0.89	J	1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Ethylbenzene	Target	810		ug/L	810	D	100.0	YES	S4VEM
o-xylene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
m,p-Xylene	Target	2600		ug/L	2600	D	100.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	6.5		ug/L	6.5		1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Total Alkanes	TIC	3.5	BN	ug/L	3.5	BN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Benzene, 1,2,4-trimethyl-	TIC	11	JN	ug/L	11	JN	1.0	YES	NV
Benzene, 1,2,3-trimethyl-	TIC	23	JN	ug/L	23	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG8	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location: A11-MW002	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 12:10:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	4.4	J	ug/L	4.4	J	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	2.7	J	ug/L	2.7	J	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	11		ug/L	11		1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.4		ug/L	6.4		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	32		ug/L	32		1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	1.0	J	ug/L	1.0	J	1.0	YES	S4VEM
Cyclohexane	Target	120		ug/L	120		1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Target	0.86	J	ug/L	0.86	J	1.0	YES	S4VEM
Methylcyclohexane	Target	570	J	ug/L	570	JD	500.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Toluene	Target	68000	J	ug/L	68000	D	500.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Ethylbenzene	Target	6400		ug/L	6400	D	500.0	YES	S4VEM
o-xylene	Target	4900		ug/L	4900	D	500.0	YES	S4VEM
m,p-Xylene	Target	20000		ug/L	20000	D	500.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	98		ug/L	98		1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	5.0		ug/L	5.0		1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene, 1-ethenyl-3-ethyl-	TIC	48	JN	ug/L	48	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Benzene, tert-butyl-	TIC	2.6	JN	ug/L	2.6	JN	1.0	YES	NV
Benzene, 2-ethyl-1,4-dimethyl-	TIC	28	JN	ug/L	28	JN	1.0	YES	NV
3-Hexanol, 2,3-dimethyl-	TIC	8.8	JN	ug/L	8.8	JN	1.0	YES	NV
Benzene, 1,2,4,5-tetramethyl-	TIC	39	JN	ug/L	39	JN	1.0	YES	NV
1H-Indene, 2,3-dihydro-4-methyl-	TIC	9.3	JN	ug/L	9.3	JN	1.0	YES	NV
Naphthalene, 1,2,3,4-tetrahydro-	TIC	43	JN	ug/L	43	JN	1.0	YES	NV
Benzene, 1-methyl-3-(1-methylethyl)	TIC	33	JN	ug/L	33	JN	1.0	YES	NV
Azulene	TIC	43	JN	ug/L	43	JN	1.0	YES	NV
Benzene, 2-propenyl-	TIC	54	JN	ug/L	54	JN	1.0	YES	NV
Benzene, 1-ethyl-2,3-dimethyl-	TIC	68	JN	ug/L	68	JN	1.0	YES	NV
2-Hexyne, 4-methyl-	TIC	3.6	JN	ug/L	3.6	JN	1.0	YES	NV
Benzeneacetaldehyde, .alpha.-methyl	TIC	19	JN	ug/L	19	JN	1.0	YES	NV
Benzene, 1,2,4-trimethyl-	TIC	180	JN	ug/L	180	JN	1.0	YES	NV
n-Butyl ether	TIC	24	JN	ug/L	24	JN	1.0	YES	NV
Benzene, 1,2-diethyl-	TIC	10	JN	ug/L	10	JN	1.0	YES	NV
Benzene, 1,2,3-trimethyl-	TIC	130	JN	ug/L	130	JN	1.0	YES	NV
o-Cymene	TIC	44	JN	ug/L	44	JN	1.0	YES	NV
Benzene, (2-methylpropyl)-	TIC	4.5	JN	ug/L	4.5	JN	1.0	YES	NV
Total Alkanes	TIC	520	BN	ug/L	520	BN	1.0	YES	NV
p-Cymene	TIC	9.7	JN	ug/L	9.7	JN	1.0	YES	NV
Benzene, propyl-	TIC	95	JN	ug/L	95	JN	1.0	YES	NV
Mesitylene	TIC	390	JN	ug/L	390	JN	1.0	YES	NV
4-Heptanone, 2,6-dimethyl-	TIC	49	JN	ug/L	49	JN	1.0	YES	NV
Benzene, 1,2,3,4-tetramethyl-	TIC	22	JN	ug/L	22	JN	1.0	YES	NV
Benzene, 1-ethyl-2-methyl-	TIC	110	JN	ug/L	110	JN	1.0	YES	NV
Benzene, 1-ethyl-3-methyl-	TIC	310	JN	ug/L	310	JN	1.0	YES	NV
Benzene, 1-methyl-4-propyl-	TIC	25	JN	ug/L	25	JN	1.0	YES	NV
Pentalene, octahydro-, cis-	TIC	16	JN	ug/L	16	JN	1.0	YES	NV
1H-Indene, octahydro-, cis-	TIC	4.9	JN	ug/L	4.9	JN	1.0	YES	NV
2-Heptanone, 4,6-dimethyl-	TIC	6.2	JN	ug/L	6.2	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG8MS	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location:	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 12:10:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	4.2	J	ug/L	4.2	J	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Spike	50		ug/L	50		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	2.8	J	ug/L	2.8	J	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	19		ug/L	19		1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.1		ug/L	6.1		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	30		ug/L	30		1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	0.97	J	ug/L	0.97	J	1.0	YES	S4VEM
Cyclohexane	Target	100		ug/L	100		1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Spike	46		ug/L	46		1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Spike	48		ug/L	48		1.0	YES	S4VEM
Methylcyclohexane	Target	590	J	ug/L	590	E	1.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Toluene	Spike	7200	J	ug/L	7200	E	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Spike	50		ug/L	50		1.0	YES	S4VEM
Ethylbenzene	Target	2500	J	ug/L	2500	E	1.0	YES	S4VEM
o-xylene	Target	4000	J	ug/L	4000	E	1.0	YES	S4VEM
m,p-Xylene	Target	8700	J	ug/L	8700	E	1.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	95		ug/L	95		1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	4.9	J	ug/L	4.9	J	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG8MSD	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location:	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 12:10:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	4.2	J	ug/L	4.2	J	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Spike	52		ug/L	52		1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	3.0	J	ug/L	3.0	J	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	21		ug/L	21		1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	6.3		ug/L	6.3		1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	32		ug/L	32		1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	1.0	J	ug/L	1.0	J	1.0	YES	S4VEM
Cyclohexane	Target	110		ug/L	110		1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Spike	47		ug/L	47		1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Spike	49		ug/L	49		1.0	YES	S4VEM
Methylcyclohexane	Target	630	J	ug/L	630	E	1.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Toluene	Spike	7500	J	ug/L	7500	E	1.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Spike	52		ug/L	52		1.0	YES	S4VEM
Ethylbenzene	Target	2600	J	ug/L	2600	E	1.0	YES	S4VEM
o-xylene	Target	4300	J	ug/L	4300	E	1.0	YES	S4VEM
m,p-Xylene	Target	8900	J	ug/L	8900	E	1.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	98		ug/L	98		1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	4.6	J	ug/L	4.6	J	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Total Alkanes	TIC		N	ug/L		N	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Sample Number: E3YG9	Method: Volatile Organics	Matrix: Water	MA Number:
Sample Location: A11-MW004A	pH: 1.0	Sample Date: 06/10/2020	Sample Time: 14:25:00
% Moisture:		% Solids: 0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Dichlorodifluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Vinyl chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromomethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichlorofluoromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloro-1,2,2-trifluoroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Acetone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Carbon disulfide	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl Acetate	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methylene chloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
trans-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Methyl tert-butyl Ether	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,2-Dichloroethene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
2-Butanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Bromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chloroform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,1-Trichloroethane	Target	3.7	J	ug/L	3.7	J	1.0	YES	S4VEM
Cyclohexane	Target	0.83	J	ug/L	0.83	J	1.0	YES	S4VEM
Carbon tetrachloride	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Trichloroethene	Target	1.0	J	ug/L	1.0	J	1.0	YES	S4VEM
Methylcyclohexane	Target	21		ug/L	21		1.0	YES	S4VEM
1,2-Dichloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromodichloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
cis-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
4-Methyl-2-pentanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Toluene	Target	52000		ug/L	52000	D	800.0	YES	S4VEM
trans-1,3-Dichloropropene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,1,2-Trichloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Tetrachloroethene	Target	5.3		ug/L	5.3		1.0	YES	S4VEM
2-Hexanone	Target	10	U	ug/L	10	U	1.0	YES	S4VEM
Dibromochloromethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromoethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Chlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Ethylbenzene	Target	330	J	ug/L	330	JD	200.0	YES	S4VEM
o-xylene	Target	71		ug/L	71		1.0	YES	S4VEM
m,p-Xylene	Target	460	J	ug/L	460	JD	200.0	YES	S4VEM
Styrene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Bromoform	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Isopropylbenzene	Target	2.1	J	ug/L	2.1	J	1.0	YES	S4VEM
1,1,2,2-Tetrachloroethane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,3-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,4-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2-Dibromo-3-chloropropane	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,4-trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
1,2,3-Trichlorobenzene	Target	5.0	U	ug/L	5.0	U	1.0	YES	S4VEM
Benzene, 1,2,4,5-tetramethyl-	TIC	3.5	JN	ug/L	3.5	JN	1.0	YES	NV



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 48947/EPW14030/E3YF9

Lab Name: Chemtech Consulting Group

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
Benzene, 1,2,3-trimethyl-	TIC	9.6	JN	ug/L	9.6	JN	1.0	YES	NV
Benzene, 1-ethyl-3-methyl-	TIC	4.9	JN	ug/L	4.9	JN	1.0	YES	NV
Benzene, 1-ethyl-2-methyl-	TIC	3.1	JN	ug/L	3.1	JN	1.0	YES	NV
Total Alkanes	TIC	110	BN	ug/L	110	BN	1.0	YES	NV
Benzene, propyl-	TIC	3.6	JN	ug/L	3.6	JN	1.0	YES	NV
Mesitylene	TIC	3.3	JN	ug/L	3.3	JN	1.0	YES	NV
Benzene, 4-ethyl-1,2-dimethyl-	TIC	3.3	JN	ug/L	3.3	JN	1.0	YES	NV



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 20060290  
**Laboratory:** STAT Analysis Corporation / Eurofins Test America  
**Matrix:** Groundwater  
**Collection date:** 06/09/20  
**Analysis/Methods:**

Wet Chemistry:  
 Nitrogen, Nitrate EPA 300.0  
 Sulfate EPA 300.0  
 Alkalinity M2320 B  
 Dissolved Gases - Methane - RSK-175

**Samples in SDG:**

<u>STAT Lab ID</u>	<u>Sample Number</u>
20060290-001	A11-MW001-200609
20060290-002	A11-MW004B-200609
20060290-003	A11-MW005-200609
20060290-004	A11-MW006-200609

Data validation was performed in accordance with the specific analytical methods, National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017), and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?			N/A
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)		Yes	
Laboratory Control Spike Duplicates RPD within limits?		N/A	
Laboratory Duplicate RPDs within limits?		N/A	
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b>Sample</b>	<b>Duplicate</b>	<b>%RPD</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
N/A					

<b>MS/MSD</b>	<b>%RPD</b>	<b>Limit</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
Nitrogen, Nitrate 20060338-003BMS/BMSD	Acceptable	20%		
Sulfate 20060338-003BMS/BMSD	Acceptable	20%		
Alkalinity 20060290-003BMS/BMSD	Acceptable	20%		

<b>LCS/LCSD</b>	<b>%RPD</b>	<b>Limits</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
N/A				

<b>Laboratory Duplicate</b>	<b>%RPD</b>	<b>Limits</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
N/A				

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	No		
Were the Field Blanks results all < RL?	N/A		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			



<b>Blanks</b>		<u>Concentration</u>	<u>MDL / RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate ICMBW1 061120	Nitrogen	0.052 J	0.2	None	Sample results nondetect or > RL
Sulfate ICMBW1 061120	Sulfate	0.377 J	4.0	None	Sample results > RL
Alkalinity ALKMBW1 061420		Nondetect		None	Sample results > RL
<b>ICB/CCB</b>		<u>Concentration</u>	<u>MDL / RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
ICB	Nitrogen, Nitrate	0.058	0.2	None	Sample results nondetect or > RL
ICB	Sulfate	0.363	4.0	None	Sample results > RL
CCB	Nitrogen, Nitrate	0.055	0.2	None	Sample results nondetect or > RL
CCB	Sulfate	0.354	4.0	None	Sample results > RL
<b>Field Blank</b>		<u>Concentration</u>	<u>MDL / RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<b>Surrogates</b>		<u>%R</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<b>MS/MSD</b>		<u>%R</u>	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate 20060338-003BMS/BMSD		Acceptable	90-110		
Sulfate 20060338-003BMS/BMSD		Acceptable	90-110		
Alkalinity 20060290-003BMS/BMSD		Acceptable	75-125		
<b>LCS/LCSD</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate ICLCSW1 061120		Acceptable	90-110		
Sulfate ICLCSW1 061120		Acceptable	90-110		
Alkalinity ALKLCSW1 061420		Acceptable	80-120		
<b>ICV</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
6/11/2020 9:14	Nitrogen, Nitrate Sulfate	73.04 Acceptable	90-110	J / UJ	All samples
<b>CCV</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
6/11/2020 '11:52	Nitrogen, Nitrate Sulfate	76.88 Acceptable	90-110	J / UJ	All samples
<b>Tune</b>					
N/A					
<b>Internal Standards</b>		<u>Area</u>	<u>Area Lower / Upper Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					



**Methane (RSK-175)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?			N/A
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)		Yes	
Laboratory Control Spike Duplicates RPD within limits?		Yes	
Laboratory Duplicate RPDs within limits?			N/A
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b><u>Sample</u></b>	<b><u>Duplicate</u></b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A					
<b>MS/MSD</b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>		<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
680-184999-3 MS / MSD (20060290-003)	Acceptable				
<b>LCS/LCSD</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>		<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Methane					
LCS 680-623376/ 3 / 4	Acceptable				
LCS 680-623376/ 6 / 7	Acceptable				
<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>		<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A					

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)		Yes	
Laboratory Control Sample criteria met?		Yes	
Were the Laboratory Method Blank results all < RL?		Yes	
Were the Field Blanks results all < RL?		N/A	
Was the ICAL criteria met?		Yes	
Was the CCV criteria met?		Yes	
Was the Tuning criteria met?		N/A	
Were the Surrogate % recoveries within laboratory determined control limits?		N/A	
Were the Internal Standard areas within ± 50 - 150%?		N/A	
<u>Comments (note deviations):</u>			

<b>Blanks</b>	<b><u>Concentration</u></b> <b><u>(mg/L)</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Methane				
MB 680-623376/ 8	Nondetect			
<b>Field Blank</b>	<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				
<b>Surrogates</b>	<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				
<b>MS/MSD</b>	<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
680-184999-3 MS / MSD (20060290-003)	Acceptable			
<b>LCS/LCSD</b>	<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Methane				
LCS 680-623376/ 3 / 4	Acceptable			
LCS 680-623376/ 6 / 7	Acceptable			
<b>ICAL</b>	<b><u>RRF</u></b>	<b><u>%RSD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
2/17/2020 8:45	Acceptable	Acceptable		
3/04/2020 9:12	Acceptable	Acceptable		



CCV	RRF	%D	Limits	Qualifiers	Associated Samples
6/22/2020 15:34	Acceptable	Acceptable			
6/22/2020 15:47	Acceptable	Acceptable			
6/22/2020 18:44	Acceptable	Acceptable			
6/22/2020 18:57	Acceptable	Acceptable			

Tune  
N/A

Internal Standards	Area	Area Lower / Upper Limit	Qualifiers	Associated Samples
N/A				

#### Representativeness:

Were sampling procedures and design criteria met?

Were holding times met?

Was preservation criteria met? (0° C - 6° C)

Were Chain-of-Custody records complete and provided in data package?

Comments (note deviations): The cooler temperatures were 4.6 and 2.1° C.

Yes No N/A

Yes

Yes

Yes

Yes

Preservation	Cooler Temperature (Degrees C)	Preservation Criteria	Qualifier	Associated Samples
	Acceptable			

Holding Times	Analyte	Days to Extraction	HT Criteria	Qualifier	Associated Samples
		Acceptable			

#### Comparability:

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

Yes No N/A

Yes

#### Completeness (90%):

Are all data in this SDG usable?

Comments (note deviations):

Yes No N/A

Yes

#### Sensitivity:

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):

Yes No N/A

Yes

Yes

#### Comment:

Data is usable with appropriate qualifiers applied.

Data Validator:

*Kristine Molloy*

Date: 1/8/2021

Data Reviewer:

Cherie Zakowski

Date: 1/12/2021



**STAT Analysis Corporation**

2242 West Harrison St., Suite 200, Chicago, IL 60612-3766

Tel: (312) 733-0551 Fax: (312) 733-2386 STATinfo@STATAnalysis.com

Accreditations: IEPA ELAP 100445; ORELAP IL300001; AIHA-LAP, LLC 101160; NVLAP LabCode 101202-0

Date Reported: July 06, 2020

Date Printed: July 06, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Quarterly GW Sampling, Work Order: 20060290 Revision 0

Lab ID: 20060290-001

Collection Date: 6/9/2020 8:40:00 AM

Client Sample ID A11-MW001-200609

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	1.8	0.20	*	mg/L	1	6/11/2020
Sulfate	25	4.0	*	mg/L	1	6/11/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	360	20		mg/L CaCO <sub>3</sub>	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	ND	0.00058		mg/L	1	6/22/2020

Lab ID: 20060290-002

Collection Date: 6/9/2020 1:05:00 PM

Client Sample ID A11-MW004B-200609

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	1.1	0.20	*	mg/L	1	6/11/2020
Sulfate	19	4.0	*	mg/L	1	6/11/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	340	20		mg/L CaCO <sub>3</sub>	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	0.035	0.00058		mg/L	1	6/22/2020

Lab ID: 20060290-003

Collection Date: 6/9/2020 4:15:00 PM

Client Sample ID A11-MW005-200609

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	2.7	0.20	*	mg/L	1	6/11/2020
Sulfate	33	4.0	*	mg/L	1	6/11/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO <sub>3</sub> )	370	20		mg/L CaCO <sub>3</sub>	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	ND	0.00058		mg/L	1	6/22/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

HT - Sample received past holding time

\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H - Holding time exceeded



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Tel: (312) 733-0551 Fax: (312) 733-2386 STATinfo@STATAnalysis.com

Accreditations: IEPA ELAP 100445; ORELAP IL300001; AIHA-LAP, LLC 101160; NVLAP LabCode 101202-0

Date Reported: July 06, 2020

Date Printed: July 06, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Quarterly GW Sampling, Work Order: 20060290 Revision 0

Lab ID: 20060290-004

Collection Date: 6/9/2020 11:00:00 AM

Client Sample ID A11-MW006-200609

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: <b>6/11/2020</b> Analyst: <b>CAB</b>
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	6/11/2020
Sulfate	7.1	4.0	*	mg/L	1	6/11/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: <b>6/14/2020</b> Analyst: <b>MD</b>
Alkalinity, Total (As CaCO3)	460	20		mg/L CaCO3	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: <b>SUB</b>
Methane	3.8	0.39		mg/L	1	6/22/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
HT - Sample received past holding time  
\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H - Holding time exceeded







<b>Blanks</b>		<u>Concentration</u>	<u>MDL / RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate ICMBW1 061120	Nitrogen	0.052 J	0.2		Sample results nondetect or > RL
Sulfate ICMBW1 061120	Sulfate	0.377 J	4.0	None	Sample results nondetect or > RL
Alkalinity ALKMBW1 061420		Nondetect			Sample results > RL
<b>ICB/CCB</b>		<u>Concentration</u>	<u>MDL / RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
ICB	Nitrogen, Nitrate	0.058	0.2	None	Sample results nondetect or > RL
ICB	Sulfate	0.363	4.0	None	
CCB	Nitrogen, Nitrate	0.072	0.2	None	Sample results nondetect or > RL
CCB	Sulfate	0.358	4.0	None	
CCB	Nitrogen, Nitrate	0.076	0.2	None	Sample results nondetect or > RL
CCB	Sulfate	0.361	4.0	None	
<b>Field Blank</b>		<u>Concentration</u>	<u>MDL / RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<b>Surrogates</b>		<u>%R</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<b>MS/MSD</b>		<u>%R</u>	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate 20060338-003BMS/BMSD		Acceptable	90-110		
Sulfate 20060338-003BMS/BMSD		Acceptable	90-110		
Alkalinity 20060290-003BMS/BMSD		Acceptable	75-125		
<b>LCS/LCSD</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Nitrogen, Nitrate ICLCSW1 061120		Acceptable	90-110		
Sulfate ICLCSW1 061120		Acceptable	90-110		
Alkalinity ALKLCSW1 061420		Acceptable	80-120		
<b>ICV</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
6/11/2020 9:14	Nitrogen, Nitrate Sulfate	73.04 Acceptable	90-110	J / UJ	All samples
<b>CCV</b>		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
6/12/2020 '2:31	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			
6/12/2020 '5:09	Nitrogen, Nitrate Sulfate	Acceptable Acceptable			



Tune  
N/A

Internal Standards  
N/A

Area

Area Lower / Upper  
Limit

Qualifiers Associated Samples



**Methane (RSK-175)**

					<u>Yes</u>	<u>No</u>	<u>N/A</u>
<b>Precision:</b>							
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?					Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)					N/A		
Laboratory Control Spike Duplicates RPD within limits?					Yes		
Laboratory Duplicate RPDs within limits?					N/A		
<u>Comments (note deviations):</u>							

<u>Field Duplicates</u>	<u>Sample</u>	<u>Duplicate</u>	<u>%RPD</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
	A11-MW007-200610	A11-MW007-200610-D	Acceptable		
<u>MS/MSD</u>	<u>%RPD</u>	<u>Limit</u>		<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					
<u>LCS/LCSD</u>	<u>%RPD</u>	<u>Limits</u>		<u>Qualifiers</u>	<u>Associated Samples</u>
Methane					
LCS 680-623522/ 3 / 4	Acceptable				
LCS 680-623522/ 6 / 7	Acceptable				
<u>Laboratory Duplicate</u>	<u>%RPD</u>	<u>Limits</u>		<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					

					<u>Yes</u>	<u>No</u>	<u>N/A</u>
<b>Accuracy:</b>							
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)						N/A	
Laboratory Control Sample criteria met?					Yes		
Were the Laboratory Method Blank results all < RL?					Yes		
Were the Field Blanks results all < RL?					N/A		
Was the ICAL criteria met?					Yes		
Was the CCV criteria met?					Yes		
Was the Tuning criteria met?					N/A		
Were the Surrogate % recoveries within laboratory determined control limits?					N/A		
Were the Internal Standard areas within ± 50 - 150%?					N/A		
<u>Comments (note deviations):</u>							

<u>Blanks</u>	<u>Concentration (mg/L)</u>	<u>MDL /RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Methane				
MB 680-623522/ 8	Nondetect			
<u>Field Blank</u>	<u>Concentration</u>	<u>MDL /RL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				
<u>Surrogates</u>	<u>%R</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				
<u>MS/MSD</u>	<u>%R</u>	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				
<u>LCS/LCSD</u>	<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
Methane				
LCS 680-623522/ 3 / 4	Acceptable			
LCS 680-623522/ 6 / 7	Acceptable			
<u>ICAL</u>	<u>RRF</u>	<u>%RSD</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
2/17/2020 8:45	Acceptable	Acceptable		
3/04/2020 9:12	Acceptable	Acceptable		



<u>CCV</u>	<u>RRF</u>	<u>%D</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
6/23/2020 13:44	Acceptable	Acceptable			
6/23/2020 14:35	Acceptable	Acceptable			
6/23/2020 17:25	Acceptable	Acceptable			
6/23/2020 17:38	Acceptable	Acceptable			

<b>Tune</b>
N/A

<u>Internal Standards</u>	<u>Area</u>	<u>Area Lower / Upper Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				

<b>Representativeness:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Were sampling procedures and design criteria met?	Yes		
Were holding times met?	Yes		
Was preservation criteria met? (0° C - 6° C)	Yes		
Were Chain-of-Custody records complete and provided in data package?	Yes		
<u>Comments (note deviations):</u> The cooler temperatures were 2.5 and 5.2° C.			

<u>Preservation</u>	<u>Cooler Temperature (Degrees C)</u>	<u>Preservation Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
	Acceptable			

<u>Holding Times</u>	<u>Analyte</u>	<u>Days to Extraction</u>	<u>HT Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
		Acceptable			

<b>Comparability:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?	Yes		
<u>Comments (note deviations):</u>			

<b>Completeness (90%):</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are all data in this SDG usable?	Yes		
<u>Comments (note deviations):</u>			

<b>Sensitivity:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are MDLs present and reported?	Yes		
Do the reporting limits meet project requirements?	Yes		
<u>Comments (note deviations):</u>			

<b>Comment:</b>
Data is usable with appropriate qualifiers applied.

Data Validator: Kristine Molloy Date: 1/6/2021

Data Reviewer: Cherie Zakowski Date: 1/8/2021



**STAT Analysis Corporation**

2242 West Harrison St., Suite 200, Chicago, IL 60612-3766

Tel: (312) 733-0551 Fax: (312) 733-2386 STATinfo@STATAnalysis.com

Accreditations: IEPA ELAP 100445; ORELAP IL300001; AIHA-LAP, LLC 101160; NVLAP LabCode 101202-0

Date Reported: July 06, 2020

Date Printed: July 06, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Quarterly GW Sampling, Work Order: 20060338 Revision 0

Lab ID: 20060338-001

Collection Date: 6/10/2020 12:10:00 PM

Client Sample ID A11-MW002-200610

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	6/12/2020
Sulfate	ND	4.0	*	mg/L	1	6/12/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO3)	400	20		mg/L CaCO3	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	19	0.39		mg/L	1	6/23/2020

Lab ID: 20060338-002

Collection Date: 6/10/2020 8:05:00 AM

Client Sample ID A11-MW003-200610

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	6/12/2020
Sulfate	9.0	4.0	*	mg/L	1	6/12/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO3)	380	20		mg/L CaCO3	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	6.7	0.39		mg/L	1	6/23/2020

Lab ID: 20060338-003

Collection Date: 6/10/2020 2:25:00 PM

Client Sample ID A11-MW004A-200610

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	0.65	0.20	*	mg/L	1	6/12/2020
Sulfate	35	4.0	*	mg/L	1	6/12/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO3)	350	20		mg/L CaCO3	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	0.19	0.00058		mg/L	1	6/23/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

HT - Sample received past holding time

\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H - Holding time exceeded



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Tel: (312) 733-0551 Fax: (312) 733-2386 STATinfo@STATAnalysis.com

Accreditations: IEPA ELAP 100445; ORELAP IL300001; AIHA-LAP, LLC 101160; NVLAP LabCode 101202-0

Date Reported: July 06, 2020

Date Printed: July 06, 2020

**ANALYTICAL RESULTS**

Client: CDM Smith Inc.

Project: 239446, SE Rockford Area 11 Quarterly GW Sampling, Work Order: 20060338 Revision 0

Lab ID: 20060338-004

Collection Date: 6/10/2020 9:50:00 AM

Client Sample ID A11-MW007-200610

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	6/12/2020
Sulfate	29	4.0	*	mg/L	1	6/12/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO3)	370	20		mg/L CaCO3	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	3.8	0.39		mg/L	1	6/23/2020

Lab ID: 20060338-005

Collection Date: 6/10/2020 9:50:00 AM

Client Sample ID A11-MW007-200610-D

Matrix: Aqueous

Analyses	Result	RL	Qualifier	Units	DF	Date Analyzed
<b>Anions by Ion Chromatography</b>	<b>E300.0</b>					Prep Date: 6/11/2020 Analyst: CAB
Nitrogen, Nitrate (As N)	ND	0.20	*	mg/L	1	6/12/2020
Sulfate	28	4.0	*	mg/L	1	6/12/2020
<b>Alkalinity</b>	<b>M2320 B</b>					Prep Date: 6/14/2020 Analyst: MD
Alkalinity, Total (As CaCO3)	360	20		mg/L CaCO3	1	6/14/2020
<b>Dissolved Gases in Water</b>	<b>RSKSOP-175</b>					Prep Date: Analyst: SUB
Methane	3.9	0.39		mg/L	1	6/23/2020

**Qualifiers:**

ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
HT - Sample received past holding time  
\* - Non-accredited parameter

RL - Reporting / Quantitation Limit for the analysis  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H - Holding time exceeded



## **September 2020 Data Validation Reports and Data Packages**



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 2009006\_\_2009007

**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:** Groundwater

**Collection date:** 9/9/2020 & 9/10/2020

**Analysis/Methods:** Wet Chemistry:  
Alkalinity M2320 B

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
2009006-08	A11-FB001-200909	2009007-01	A11-MW002-200910
2009006-09	A11-MW001-200909	2009007-02	A11-MW007-200910
2009006-10	A11-MW004B-200909	2009007-03	A11-MW004A-200910
2009006-11	A11-MW006-200909	2009007-04	A11-MW007-200910-D
2009006-12	A11-MW005-200909	2009007-05	A11-MW003-200910-D
2009006-13	A11-MW130A-200909		

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters (Alkalinity 2320B)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	N/A		
Laboratory Control Spike Duplicates RPD within limits?	N/A		
Laboratory Duplicate RPDs within limits?	Yes		
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b><u>Sample</u></b>	<b><u>Duplicate</u></b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	A11-MW007-200910	A11-MW007-200910-D	Acceptable		

<b>MS/MSD</b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>LCS/LCSD</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20I015-DUP1	Acceptable			

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	N/A		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	Yes		
Was the ICAL criteria met?	N/A		
Was the CCV criteria met?	N/A		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

<b>Blanks</b>	<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20I015-BLK1	Nondetect			



<b>Field Blank</b> A11-FB001-200909	<u>Concentration</u> Nondetect	<u>MDL / PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>Surrogates</b> N/A	<u>%R</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>MS/MSD</b> N/A	<u>%R</u>	<u>Limits (%)</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>LCS/LCSD</b> B20I015-SRM1	<u>%R</u> Acceptable	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
<b>ICV</b> N/A		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u> <u>Associated Samples</u>
<b>CCV</b> N/A		<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u> <u>Associated Samples</u>
<b>Tune</b> N/A				
<b>Internal Standards</b> N/A	<u>Area</u>	<u>Area Lower / Upper</u> <u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>

**Representativeness:**

	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Were sampling procedures and design criteria met?	Yes		
Were holding times met?	Yes		
Was preservation criteria met? (0° C - 6° C)	Yes		
Were Chain-of-Custody records complete and provided in data package?	Yes		
<u>Comments (note deviations):</u> The cooler temperatures were 1.1 and 3.8° C.			

<b>Preservation</b>	<u>Cooler</u> <u>Temperature</u> <u>(Degrees C)</u> Acceptable	<u>Preservation</u> <u>Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
<b>Holding Times</b>	<u>Analyte</u>	<u>Days to Extraction</u> Acceptable	<u>HT Criteria</u>	<u>Qualifier</u> <u>Associated Samples</u>

**Comparability:**

	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?	Yes		
<u>Comments (note deviations):</u>			

**Completeness (90%):**

	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Are all data in this SDG usable?	Yes		
<u>Comments (note deviations):</u>			

**Sensitivity:**

	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Are MDLs present and reported?	Yes		
Do the reporting limits meet project requirements?	Yes		
<u>Comments (note deviations):</u>			

**Comment:**

Data is usable as reported.

Data Validator:

*Kristine Molloy*

Date: 1/22/2021

Data Reviewer:

Cherie Zakowski

Date: 1/25/2021





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Oct-19-20 10:41

**Alkalinity by SM 2320B**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-FB001-200909 (2009006-08)**

**Matrix: Water**

**Sampled: Sep-09-20 18:00**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>U</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW001-200909 (2009006-09)**

**Matrix: Water**

**Sampled: Sep-09-20 16:25**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>340</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW004B-200909 (2009006-10)**

**Matrix: Water**

**Sampled: Sep-09-20 16:30**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>340</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW006-200909 (2009006-11)**

**Matrix: Water**

**Sampled: Sep-09-20 11:45**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>440</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW005-200909 (2009006-12)**

**Matrix: Water**

**Sampled: Sep-09-20 13:40**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>370</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW130A-200909 (2009006-13)**

**Matrix: Water**

**Sampled: Sep-09-20 09:55**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>330</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW002-200910 (2009007-01)**

**Matrix: Water**

**Sampled: Sep-10-20 13:05**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>420</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW007-200910 (2009007-02)**

**Matrix: Water**

**Sampled: Sep-10-20 10:55**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>530</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Oct-19-20 10:41

**Alkalinity by SM 2320B**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW004A-200910 (2009007-03)**

**Matrix: Water**

**Sampled: Sep-10-20 15:50**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>340</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW007-200910-D (2009007-04)**

**Matrix: Water**

**Sampled: Sep-10-20 10:55**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>530</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20

**A11-MW003-200910-D (2009007-05)**

**Matrix: Water**

**Sampled: Sep-10-20 08:45**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>370</b>			20	mg CaCO3/L	1	B20I015	Sep-15-20	Sep-15-20





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Oct-19-20 10:41

**Notes and Definitions**

- \* This Quality Control measure meets the requirements of the CRL SOP for this analyte.
- U Not Detected
- NR Not Reported
- Q QC limit Exceeded



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 680-188662

**Laboratory:** Eurofins Test America

**Matrix:** Groundwater

**Collection date:** 09/09/2020 & 09/10/2020

**Analysis/Methods:** Dissolved Gases - Methane - RSK-175

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
680-188662-1	A11-MW006-200909	680-188662-7	A11-MW003-200910
680-188662-2	A11-MW130A-200909	680-188662-8	A11-MW007-200910
680-188662-3	A11-MW005-200909	680-188662-9	A11-MW007-200910-D
680-188662-4	A11-MW001-200909	680-188662-10	A11-MW002-200910
680-188662-5	A11-MW004B-200909	680-188662-11	A11-MW004A-200910
680-188662-6	A11-FB01-200909	680-188662-12	A11-TB001-200909

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Methane (RSK-175)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	N/A		
Laboratory Control Spike Duplicates RPD within limits?	Yes		
Laboratory Duplicate RPDs within limits?	N/A		
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b><u>Sample</u></b>	<b><u>Duplicate</u></b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	A11-MW007-200910	A11-MW007-200910-D	Acceptable		

<b>MS/MSD</b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>LCS/LCSD</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
LCS 680-635562/ 3 / 4	Acceptable			
LCS 680-635562/ 6 / 7	Acceptable			

<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	No		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

<b>Blanks</b>	<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
MB 680-635562/ 8	(mg/L) Nondetect			



<b>Field Blank</b>		<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
A11-FB01-200909	Methane	0.62	0.29 / 0.58	U-RL	680-188662-2		
A11-TB001-200909	Methane	0.64	0.29 / 0.58	U-RL	680-188662-2		
<b>Surrogates</b>		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
N/A							
<b>MS/MSD</b>		<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
N/A							
<b>LCS/LCSD</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
LCS 680-635562/ 3 / 4		Acceptable					
LCS 680-635562/ 6 / 7		Acceptable					
<b>ICAL</b>		<b><u>RRF</u></b>	<b><u>%RSD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
2/17/2020 8:45		Acceptable	Acceptable				
3/04/2020 9:12		Acceptable	Acceptable				
<b>ICV / CCV</b>		<b><u>RRF</u></b>	<b><u>%D</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
ICV							
3/04/2020 11:29		Acceptable	Acceptable				
CCV							
09/23/2020 10:40		Acceptable	Acceptable				
09/23/2020 11:19		Acceptable	Acceptable				
09/23/2020 15:15		Acceptable	Acceptable				
09/23/2020 15:28		Acceptable	Acceptable				
09/23/2020 17:19		Acceptable	Acceptable				
09/23/2020 17:32		Acceptable	Acceptable				
<b>Tune</b>							
N/A							
<b>Internal Standards</b>		<b><u>Area</u></b>	<b><u>Area Lower / Upper Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
N/A							
<b>Representativeness:</b>					<b><u>Yes</u></b>	<b><u>No</u></b>	<b><u>N/A</u></b>
Were sampling procedures and design criteria met?						Yes	
Were holding times met?						Yes	
Was preservation criteria met? (0° C - 6° C)						No	
Were Chain-of-Custody records complete and provided in data package?						Yes	
<u>Comments (note deviations):</u> The cooler temperature was 17.9° C.							
<b>Preservation</b>		<b><u>Cooler Temperature (Degrees C)</u></b>	<b><u>Preservation Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>		
	Methane	17.9	0 - 6 ° C	J /UJ	All samples		
<b>Holding Times</b>	<b><u>Analyte</u></b>	<b><u>Days to Extraction</u></b>	<b><u>HT Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>		
		Acceptable					
<b>Comparability:</b>					<b><u>Yes</u></b>	<b><u>No</u></b>	<b><u>N/A</u></b>
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?						Yes	
<u>Comments (note deviations):</u>							
<b>Completeness (90%):</b>					<b><u>Yes</u></b>	<b><u>No</u></b>	<b><u>N/A</u></b>
Are all data in this SDG usable?						Yes	
<u>Comments (note deviations):</u>							



Sensitivity:	Yes	No	N/A
Are MDLs present and reported?	Yes		
Do the reporting limits meet project requirements?	Yes		

Comments (note deviations):

Comment:

As noted by the laboratory, samples were received properly preserved on ice and in good condition, however, water was present in the cooler, indicating melted ice.

Data is usable with appropriate qualifiers applied.

Data Validator:

Kristine Molloy

Date: 1/22/2021

Data Reviewer:

Cherie Zakowski

Date: 1/25/2021



# Detection Summary

Client: CDM Smith, Inc.  
Project/Site: Methane Analysis - SE Rockford Area 11

Job ID: 680-188662-1

## Client Sample ID: A11-MW006-200909

## Lab Sample ID: 680-188662-1

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	4100		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW130A-200909

## Lab Sample ID: 680-188662-2

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.57	J	0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW005-200909

## Lab Sample ID: 680-188662-3

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.94		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW001-200909

## Lab Sample ID: 680-188662-4

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.82		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW004B-200909

## Lab Sample ID: 680-188662-5

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	27		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-FB01-200909

## Lab Sample ID: 680-188662-6

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.62		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW003-200910

## Lab Sample ID: 680-188662-7

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	3500		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW007-200910

## Lab Sample ID: 680-188662-8

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	25000		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW007-200910-D

## Lab Sample ID: 680-188662-9

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	21000		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW002-200910

## Lab Sample ID: 680-188662-10

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	26000		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW004A-200910

## Lab Sample ID: 680-188662-11

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	160		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-TB001-200909

## Lab Sample ID: 680-188662-12

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.64		0.58	0.29	ug/L	1			RSK-175	Total/NA

This Detection Summary does not include radiochemical test results.

Eurofins TestAmerica, Savannah



# Client Sample Results

Client: CDM Smith, Inc.  
Project/Site: Methane Analysis - SE Rockford Area 11

Job ID: 680-188662-1

**Client Sample ID: A11-MW006-200909**

**Lab Sample ID: 680-188662-1**

Date Collected: 09/09/20 11:45

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	4100		390	39	ug/L			09/23/20 13:06	1

**Client Sample ID: A11-MW130A-200909**

**Lab Sample ID: 680-188662-2**

Date Collected: 09/09/20 09:55

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.57	J	0.58	0.29	ug/L			09/23/20 13:19	1

**Client Sample ID: A11-MW005-200909**

**Lab Sample ID: 680-188662-3**

Date Collected: 09/09/20 13:40

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.94		0.58	0.29	ug/L			09/23/20 13:32	1

**Client Sample ID: A11-MW001-200909**

**Lab Sample ID: 680-188662-4**

Date Collected: 09/09/20 16:25

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.82		0.58	0.29	ug/L			09/23/20 13:45	1

**Client Sample ID: A11-MW004B-200909**

**Lab Sample ID: 680-188662-5**

Date Collected: 09/09/20 16:30

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	27		0.58	0.29	ug/L			09/23/20 13:58	1

**Client Sample ID: A11-FB01-200909**

**Lab Sample ID: 680-188662-6**

Date Collected: 09/09/20 18:00

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.62		0.58	0.29	ug/L			09/23/20 14:11	1

**Client Sample ID: A11-MW003-200910**

**Lab Sample ID: 680-188662-7**

Date Collected: 09/10/20 08:45

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	3500		390	39	ug/L			09/23/20 16:01	1



# Client Sample Results

Client: CDM Smith, Inc.

Job ID: 680-188662-1

Project/Site: Methane Analysis - SE Rockford Area 11

**Client Sample ID: A11-MW007-200910**

**Lab Sample ID: 680-188662-8**

Date Collected: 09/10/20 10:55

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	25000		390	39	ug/L			09/23/20 14:24	1

**Client Sample ID: A11-MW007-200910-D**

**Lab Sample ID: 680-188662-9**

Date Collected: 09/10/20 10:55

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	21000		390	39	ug/L			09/23/20 14:37	1

**Client Sample ID: A11-MW002-200910**

**Lab Sample ID: 680-188662-10**

Date Collected: 09/10/20 13:05

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	26000		390	39	ug/L			09/23/20 14:49	1

**Client Sample ID: A11-MW004A-200910**

**Lab Sample ID: 680-188662-11**

Date Collected: 09/10/20 15:50

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	160		0.58	0.29	ug/L			09/23/20 15:02	1

**Client Sample ID: A11-TB001-200909**

**Lab Sample ID: 680-188662-12**

Date Collected: 09/09/20 08:00

Matrix: Water

Date Received: 09/14/20 08:50

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.64		0.58	0.29	ug/L			09/23/20 15:48	1



## Default Detection Limits

Client: CDM Smith, Inc.

Job ID: 680-188662-1

Project/Site: Methane Analysis - SE Rockford Area 11

### Method: RSK-175 - Dissolved Gases (GC)

Analyte	RL	MDL	Units
Methane	0.58	0.29	ug/L
Methane (TCD)	390	39	ug/L



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** E200903  
**Laboratory:** ESAT / Tech Law  
**Matrix:** Groundwater  
**Collection date:** 09/09/2020 & 09/10/2020  
**Analysis/Methods:** 1,4-Dioxane - SW-846 8000D SIM

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
E200901-01	A11-FB001-200909	E200902-01	A11-MW004A-200910
E200901-02	A11-MW001-200909	E200902-02	A11-MW007-200910-D
E200901-03	A11-MW004B-200909	E200902-03	A11-MW007-200910
E200901-04	A11-MW005-200909	E200902-04	A11-MW003-200910
E200901-05	A11-MW006-200909	E200902-05	A11-TB002-200910
E200901-06	A11-MW130A-200909	E200902-06	A11-MW002-200910
E200901-07	A11-TB001-200909		

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Volatile Organic Compounds 8260 / 1,4-Dioxane 8000D**

Precision:	Yes	No	N/A
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	Yes		
Laboratory Control Spike Duplicates RPD within limits?	Yes		
Laboratory Duplicate RPDs within limits?	N/A		
<u>Comments (note deviations):</u>			

Field Duplicates	Sample A11-MW007- 200910 ND	Duplicate A11-MW007-200910- D ND	%RPD	Qualifiers	Associated Samples
1,4-Dioxane					

MS/MSD	%RPD	Limit	Qualifiers	Associated Samples
E20I001-MS1 / MSD1 (E200901-04)	Acceptable			

LCS/LCSD	%RPD	Limits	Qualifiers	Associated Samples
E20I001-BS1 / BSD1	Acceptable			

Laboratory Duplicate	%RPD	Limits	Qualifiers	Associated Samples
N/A				

Accuracy:	Yes	No	N/A
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	No		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	Yes		
Were the Surrogate % recoveries within laboratory determined control limits?	Yes		
Were the Internal Standard areas within ± 50 - 150%?	Yes		
<u>Comments (note deviations):</u>			

Blanks	Concentration	MDL /PQL	Qualifiers	Associated Samples
E20I001-BLK1	Nondetect			



<b>Field Blank</b> A11-FB001-200909 A11-TB001-200909 A11-TB002-200910	<b>Analyte</b> 1,4-Dioxane	<b>Concentration</b> 7.81 Nondetect Nondetect	<b>MDL / PQL</b> 0.207	<b>Qualifiers</b> None	<b>Associated Samples</b> Sample results nondetect or > RL
<b>Surrogates</b>		<b>%R</b> Acceptable	<b>Limit</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
<b>MS/MSD</b> E20I001-MS1 / MSD1 (E200901-04)	1,4-Dioxane	<b>%R</b> 125 / 134	<b>Limits (%)</b> 64-112	<b>Qualifiers</b> J	<b>Associated Samples</b> E200901-04
<b>LCS/LCSD</b> E20I001-BS1 / BSD1		<b>%R</b> Acceptable	<b>Limits</b>	<b>Qualifiers</b>	<b>Associated Samples</b>
<b>ICAL</b> May 27, 2020		<b>RRF</b> Acceptable	<b>%RSD</b> Acceptable	<b>Limits</b>	<b>Qualifiers</b> <b>Associated Samples</b>
<b>ICV / CCV</b> ICV 5/27/2020 1:00  CCV 9/21/2020 11:50 9/21/2020 20:35		<b>RRF</b> Acceptable  Acceptable Acceptable	<b>%D</b> Acceptable  Acceptable Acceptable	<b>Limits</b>	<b>Qualifiers</b> <b>Associated Samples</b>
<b>Tune</b> Acceptable					
<b>MRL Check</b>  E20I001-MRL1			<b>%R</b> Acceptable	<b>Limits</b>	<b>Qualifiers</b> <b>Associated Samples</b>
<b>Internal Standards</b>	<b>Area</b>	<b>Area Lower / Upper</b> <b>Limit</b> Acceptable		<b>Qualifiers</b>	<b>Associated Samples</b>
<b>Representativeness:</b>					
Were sampling procedures and design criteria met?					<b>Yes</b> <b>No</b> <b>N/A</b>
Were holding times met?					<b>Yes</b>
Was preservation criteria met? (0° C - 6° C)					<b>Yes</b>
Were Chain-of-Custody records complete and provided in data package?					<b>Yes</b>
<u>Comments (note deviations):</u> The cooler temperatures were 1.1 & 0.9 ° C.					
<b>Preservation</b>	<b>Cooler Temperature (Degrees C)</b> Acceptable	<b>Preservation Criteria</b>		<b>Qualifier</b>	<b>Associated Samples</b>
<b>Holding Times</b>	<b>Analyte</b>	<b>Days to Extraction</b> Acceptable	<b>HT Criteria</b>	<b>Qualifier</b>	<b>Associated Samples</b>
<b>Comparability:</b>					
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?					<b>Yes</b> <b>No</b> <b>N/A</b>
<u>Comments (note deviations):</u>					



**Completeness (90%):**

Are all data in this SDG usable?

Comments (note deviations):**Yes No N/A****Yes****Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):**Yes No N/A****Yes****Yes****Comment:**

Data is usable with appropriate qualifiers applied.

Data Validator:

*Kristine Molloy*Date: 5/1/2021

Data Reviewer:

Cherie ZakowskiDate: 5/5/2021





TechLaw Inc ESAT Region 5  
536 South Clark Street, Suite 734  
Chicago, IL 60605  
(312) 353-8303  
(312) 353-5814 (Fax)  
[www.techlawinc.com](http://www.techlawinc.com)

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Michelle Kerr

Reported:  
Oct-14-20 10:13

**1,4-Dioxane by GC-MS**  
**TechLaw - ESAT Contract**

**A11-FB001-200909 (E200901-01)**

**Matrix: Water**

**Sampled: Sep-09-20 18:00**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>7.81</b>			0.207	ug/L	1	E201001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.845			81.8%		64-109	"	"	"

**A11-MW001-200909 (E200901-02)**

**Matrix: Water**

**Sampled: Sep-09-20 16:25**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.205	ug/L	1	E201001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.839			81.9%		64-109	"	"	"

**A11-MW004B-200909 (E200901-03)**

**Matrix: Water**

**Sampled: Sep-09-20 16:30**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>7.86</b>			0.207	ug/L	1	E201001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.779			75.4%		64-109	"	"	"

**A11-MW005-200909 (E200901-04)**

**Matrix: Water**

**Sampled: Sep-09-20 13:40**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>8.18</b>			0.205	ug/L	1	E201001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.801			78.2%		64-109	"	"	"

**A11-MW006-200909 (E200901-05)**

**Matrix: Water**

**Sampled: Sep-09-20 11:45**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>8.42</b>			0.203	ug/L	1	E201001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.841			82.7%		64-109	"	"	"





TechLaw Inc ESAT Region 5  
536 South Clark Street, Suite 734  
Chicago, IL 60605  
(312) 353-8303  
(312) 353-5814 (Fax)  
[www.techlawinc.com](http://www.techlawinc.com)

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Michelle Kerr

Reported:  
Oct-14-20 10:13

**1,4-Dioxane by GC-MS**  
**TechLaw - ESAT Contract**

**A11-MW130A-200909 (E200901-06)**

**Matrix: Water**

**Sampled: Sep-09-20 09:55**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>6.10</b>			0.205	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.824			80.4%		64-109	"	"	"

**A11-TB001-200909 (E200901-07)**

**Matrix: Water**

**Sampled: Sep-09-20 07:30**

**Received: Sep-10-20 10:40**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.208	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.964			92.6%		64-109	"	"	"

**A11-MW004A-200910 (E200902-01)**

**Matrix: Water**

**Sampled: Sep-10-20 15:50**

**Received: Sep-11-20 10:44**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>1.09</b>			0.203	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.737			72.5%		64-109	"	"	"

**A11-MW007-200910-D (E200902-02)**

**Matrix: Water**

**Sampled: Sep-10-20 10:55**

**Received: Sep-11-20 10:44**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.203	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.935			92.0%		64-109	"	"	"

**A11-MW007-200910 (E200902-03)**

**Matrix: Water**

**Sampled: Sep-10-20 10:55**

**Received: Sep-11-20 10:44**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.212	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.860			81.2%		64-109	"	"	"





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[www.techlawinc.com](http://www.techlawinc.com)

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Michelle Kerr

Reported:  
Oct-14-20 10:13

**1,4-Dioxane by GC-MS**  
**TechLaw - ESAT Contract**

**A11-MW003-200910 (E200902-04)**

**Matrix: Water**

**Sampled: Sep-10-20 08:45**

**Received: Sep-11-20 10:44**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>7.23</b>			0.205	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.884			86.3%		64-109	"	"	"

**A11-TB002-200910 (E200902-05)**

**Matrix: Water**

**Sampled: Sep-10-20 08:00**

**Received: Sep-11-20 10:44**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>U</b>			0.203	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.941			92.6%		64-109	"	"	"

**A11-MW002-200910 (E200902-06)**

**Matrix: Water**

**Sampled: Sep-10-20 13:05**

**Received: Sep-11-20 10:44**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>1,4-Dioxane</b>	<b>2.90</b>			0.214	ug/L	1	E20I001	Sep-18-20	Sep-21-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>1,4-Dioxane-d8</i>	0.991			92.7%		64-109	"	"	"





TechLaw Inc ESAT Region 5  
536 South Clark Street, Suite 734  
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(312) 353-8303  
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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Michelle Kerr

Reported:  
Oct-14-20 10:13

## 1,4-Dioxane by GC-MS - Quality Control

### TechLaw - ESAT Contract

#### Batch E20I001 - EPA 522

##### Blank (E20I001-BLK1)

Prepared: Sep-18-20 Analyzed: Sep-21-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	U			0.200	ug/L						
Surrogate: 1,4-Dioxane-d8	0.796				"	1.00		79.6%	64-109		

##### LCS (E20I001-BS1)

Prepared: Sep-18-20 Analyzed: Sep-21-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	0.799			0.200	ug/L	1.00		79.9%	70-106		
Surrogate: 1,4-Dioxane-d8	0.814				"	1.00		81.4%	64-109		

##### LCS Dup (E20I001-BSD1)

Prepared: Sep-18-20 Analyzed: Sep-21-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	0.794			0.200	ug/L	1.00		79.4%	70-106	0.552	17
Surrogate: 1,4-Dioxane-d8	0.807				"	1.00		80.7%	64-109		

##### MRL Check (E20I001-MRL1)

Prepared: Sep-18-20 Analyzed: Sep-21-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	0.132	J		0.200	ug/L	0.200		66.2%	49-131		
Surrogate: 1,4-Dioxane-d8	0.796				"	1.00		79.6%	64-109		

##### Matrix Spike (E20I001-MS1)

Source: E200901-04

Prepared: Sep-18-20 Analyzed: Sep-21-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	9.45	Q		0.203	ug/L	1.02	8.18	125%	64-112		
Surrogate: 1,4-Dioxane-d8	0.862				"	1.02		84.8%	64-109		

##### Matrix Spike Dup (E20I001-MSD1)

Source: E200901-04

Prepared: Sep-18-20 Analyzed: Sep-21-20

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
1,4-Dioxane	9.55	Q		0.205	ug/L	1.02	8.18	134%	64-112	6.72	12
Surrogate: 1,4-Dioxane-d8	0.895				"	1.02		87.4%	64-109		





TechLaw Inc ESAT Region 5  
536 South Clark Street, Suite 734  
Chicago, IL 60605  
(312) 353-8303  
(312) 353-5814 (Fax)  
[www.techlawinc.com](http://www.techlawinc.com)

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION  
Project Number: ILD981000417  
Project Manager: Michelle Kerr

**Reported:**  
Oct-14-20 10:13

### Notes and Definitions

J The identification of the analyte is acceptable; the reported value is an estimate.  
U Not Detected  
NR Not Reported  
Q QC limit Exceeded



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 2009006

**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:** Groundwater

**Collection date:** 09/09/20

**Analysis/Methods:** Wet Chemistry: Anions - EPA 300.0

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
2009006-08	A11-FB001-200909	2009006-11	A11-MW006-200909
2009006-09	A11-MW001-200909	2009006-12	A11-MW005-200909
2009006-10	A11-MW004B-200909	2009006-13	A11-MW130A-200909

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters (Anions 300.0)**

Precision:	Yes	No	N/A
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?			N/A
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)			N/A
Laboratory Control Spike Duplicates RPD within limits?			N/A
Laboratory Duplicate RPDs within limits?			N/A
<u>Comments (note deviations):</u>			

Field Duplicates	Sample	Duplicate	%RPD	Qualifiers	Associated Samples
N/A					

MS/MSD	%RPD	Limit	Qualifiers	Associated Samples
N/A				

LCS/LCSD	%RPD	Limits	Qualifiers	Associated Samples
N/A				

Laboratory Duplicate	%RPD	Limits	Qualifiers	Associated Samples
E20I011-DUP1	Acceptable			

Accuracy:	Yes	No	N/A
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)		No	
Laboratory Control Sample criteria met?		Yes	
Were the Laboratory Method Blank results all < RL?		Yes	
Were the Field Blanks results all < RL?		Yes	
Was the ICAL criteria met?		Yes	
Was the CCV criteria met?		Yes	
Was the Tuning criteria met?		N/A	
Were the Surrogate % recoveries within laboratory determined control limits?		N/A	
Were the Internal Standard areas within ± 50 - 150%?		N/A	
<u>Comments (note deviations):</u>			

Blanks	Concentration	MDL /PQL	Qualifiers	Associated Samples
E20I011-BLK1				
Nitrogen, Nitrate	Nondetect			
Sulfate	Nondetect			



<b>ICB/CCB</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
ICB	Nitrogen, Nitrate Sulfate	Nondetect 0.04	0.1 / 0.12	None	Sample results > RL
CCB	Nitrogen, Nitrate Sulfate	Nondetect 0.04	0.1 / 0.12	None	Sample results > RL
<b>Field Blank</b> A11-FB001-200909		<b><u>Concentration</u></b> Nondetect	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>Surrogates</b> N/A		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>MS/MSD</b> E20I011-MS1 Nitrogen, Nitrate		<b><u>%R</u></b> Acceptable	<b><u>Limits (%)</u></b> 80-120	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Sulfate		69	80-120	J- / UJ	All samples
<b>LCS/LCSD</b> E20I011-BS1 Nitrogen, Nitrate		<b><u>%R</u></b> Acceptable	<b><u>Limits</u></b> 90-110	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Sulfate		Acceptable	90-110		
<b>ICV</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Nitrogen, Nitrate Sulfate		Acceptable Acceptable			
<b>CCV</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
Nitrogen, Nitrate Sulfate		Acceptable Acceptable			
<b>MRL Check</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20I011-MRL1 Nitrogen, Nitrate Sulfate		Acceptable Acceptable			
<b>Tune</b> N/A					
<b>Internal Standards</b> N/A		<b><u>Area</u></b>	<b><u>Area Lower / Upper Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>Representativeness:</b>					<b>Yes No N/A</b>
Were sampling procedures and design criteria met?					Yes
Were holding times met?					Yes
Was preservation criteria met? (0° C - 6° C)					Yes
Were Chain-of-Custody records complete and provided in data package?					Yes
<b>Comments (note deviations):</b> The cooler temperatures were 1.1 and 3.8 ° C.					
<b>Preservation</b>		<b><u>Cooler Temperature (Degrees C)</u></b>	<b><u>Preservation Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>
		Acceptable			



Holding Times	Analyte	Days to Extraction Acceptable	HT Criteria	Qualifier	Associated Samples
<b>Comparability:</b> Were analytical procedures and methods followed as defined in the QAPP or field change documentation? <u>Comments (note deviations):</u>					<b>Yes No N/A</b> Yes
<b>Completeness (90%):</b> Are all data in this SDG usable? <u>Comments (note deviations):</u>					<b>Yes No N/A</b> Yes
<b>Sensitivity:</b> Are MDLs present and reported? Do the reporting limits meet project requirements? <u>Comments (note deviations):</u>					<b>Yes No N/A</b> Yes Yes
<b>Comment:</b> Data is usable with appropriate qualifiers applied.					
Data Validator:	<u>Kristine Molloy</u>			Date:	<u>1/20/2021</u>
Data Reviewer:	<u>Cherie Zakowski</u>			Date:	<u>1/25/2021</u>





## Environmental Protection Agency Region 5

### US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone: (312) 353-8370 Fax: (312) 886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Nov-09-20 14:18

### Anions by Ion Chromatography, EPA 300.0 (modified) US EPA Region 5 LSASD Analytical Services Branch

**A11-FB001-200909 (2009006-08)**

**Matrix: Water**

**Sampled: Sep-09-20 18:00**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	U			0.12	mg/L	1	B20I011	Sep-10-20	Sep-11-20
Nitrate - NO <sub>3</sub>	U			0.12	"	"	"	"	"

**A11-MW001-200909 (2009006-09)**

**Matrix: Water**

**Sampled: Sep-09-20 16:25**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	28.9			0.12	mg/L	1	B20I011	Sep-10-20	Sep-10-20
Nitrate - NO <sub>3</sub>	11.3			0.12	"	"	"	"	"

**A11-MW004B-200909 (2009006-10)**

**Matrix: Water**

**Sampled: Sep-09-20 16:30**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	18.8			0.12	mg/L	1	B20I011	Sep-10-20	Sep-10-20
Nitrate - NO <sub>3</sub>	5.20			0.12	"	"	"	"	"

**A11-MW006-200909 (2009006-11)**

**Matrix: Water**

**Sampled: Sep-09-20 11:45**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	5.02			0.12	mg/L	1	B20I011	Sep-10-20	Sep-10-20
Nitrate - NO <sub>3</sub>	U			0.12	"	"	"	"	"

**A11-MW005-200909 (2009006-12)**

**Matrix: Water**

**Sampled: Sep-09-20 13:40**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	25.2	(MS), L		0.12	mg/L	1	B20I011	Sep-10-20	Sep-10-20
Nitrate - NO <sub>3</sub>	9.53			0.12	"	"	"	"	"

**A11-MW130A-200909 (2009006-13)**

**Matrix: Water**

**Sampled: Sep-09-20 09:55**

**Received: Sep-10-20 10:05**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	17.5			0.12	mg/L	1	B20I011	Sep-10-20	Sep-10-20
Nitrate - NO <sub>3</sub>	5.91			0.12	"	"	"	"	"





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Nov-09-20 14:18

**Notes and Definitions**

- L The identification of the analyte is acceptable; the reported value may be biased low. The actual value is expected to be greater than the reported value.
- (MS) Matrix spike recovery criteria not met for this analyte
- U Not Detected
- NR Not Reported
- Q QC limit Exceeded



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:**  
**Laboratory:**

2009006 / 2009007

ESAT

**Matrix:**  
**Collection date:**  
**Analysis/Methods:**

Groundwater

09/09/2020 & 09/10/2020

Volatile Organic Compounds (VOCs) 8260

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
2009006-01	A11-TB001-200909	2009007-06	A11-MW007-200910
2009006-02	A11-MW004B-200909	2009007-07	A11-TB002-200910
2009006-03	A11-MW005-200909	2009007-08	A11-MW007-200910-D
2009006-04	A11-MW006-200909	2009007-09	A11-MW002-200910
2009006-05	A11-MW130A-200909	2009007-10	A11-MW004A-200910
2009006-06	A11-MW001-200909	2009007-11	A11-MW003-200910
2009006-07	A11-FB001-200909		

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Volatile Organic Compounds 8260**

Precision:	Yes	No	N/A
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	Yes		
Laboratory Control Spike Duplicates RPD within limits?	No		
Laboratory Duplicate RPDs within limits?	N/A		
<u>Comments (note deviations):</u>			

Field Duplicates	Sample A11-MW007- 200910	Duplicate A11-MW007-200910- D	%RPD	Qualifiers	Associated Samples
			Acceptable		

MS/MSD	%RPD	Limit	Qualifiers	Associated Samples
B20I010-MS1 / MSD1 (2009006-03RE1)	Acceptable			

LCS/LCSD	%RPD	Limits	Qualifiers	Associated Samples
B20I010-BS1 / BSD1	Acceptable			
B20I012-BS1/ BSD1	2,2-Dichloropropane	40.1	J**	2009006-02RE1, 2009006-04RE1 through 2009006-06RE1, 2009007-06 through 2009007-08
	Toluene	27.5	J**	
B20I014-BS1 / BSD1	2,2-Dichloropropane	68.6	J**	2009007-06RE1, 2009007-09RE1 through 2009007-11RE1, 2009007-10RE2 through 2009007- 11RE2

\*\*Qualification required for detected results only - associated results nondetect - no qualification required

Laboratory Duplicate	%RPD	Limits	Qualifiers	Associated Samples
N/A				



Accuracy:		Yes	No	N/A
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)		Yes		
Laboratory Control Sample criteria met?		No		
Were the Laboratory Method Blank results all < RL?		Yes		
Were the Field Blanks results all < RL?		Yes		
Was the ICAL criteria met?		No		
Was the CCV criteria met?		No		
Was the Tuning criteria met?		Yes		
Were the Surrogate % recoveries within laboratory determined control limits?		Yes		
Were the Internal Standard areas within ± 50 - 150%?		Yes		
Comments (note deviations):				

Blanks	Concentration	MDL /PQL	Qualifiers	Associated Samples
B20I010-BLK1	Nondetect			
B20I010-BLK2	Nondetect			
B20I012-BLK1	Nondetect			
B20I014-BLK1	Nondetect			

Field Blank	Concentration	MDL / PQL	Qualifiers	Associated Samples
A11-TB001-200909	Nondetect			
A11-FB001-200909	Nondetect			
A11-TB002-200910	Nondetect			

Surrogates	%R	Limit	Qualifiers	Associated Samples
	Acceptable			

MS/MSD	%R	Limits (%)	Qualifiers	Associated Samples
B20I010-MS1 / MSD1 (2009006-03RE1)	Acceptable			

LCS/LCSD		%R	Limits	Qualifiers	Associated Samples
B20I010-BS1 / BSD1		Acceptable			
B20I012-BS1/ BSD1	Toluene	100 / 132	70-130	J**	2009006-02RE1, 2009006-04RE1 through 2009006-06RE1, 2009007-06 through 2009007-08
B20I014-BS1 / BSD1	Dichlorodifluoromethane	59.9 / 60.1	70-130	J / UJ	2009007-06RE1, 2009007-09RE1 through 2009007-11RE1,
	2,2-Dichloropropane	99 / 48.5	70-130	J / UJ	2009007-10RE2 through 2009007-11RE2
B20I014-BS2	Dichlorodifluoromethane	64.8	70-130	J / UJ	2009007-08RE1, 2009007-09RE2
	2,2-Dichloropropane	64.8	70-130	J / UJ	

\*\*Qualification required for detected results only - associated results nondetect - no qualification required

ICAL		RRF	%RSD	Limits	Qualifiers	Associated Samples
9/10/2020	Dichlorodifluoromethane	Acceptable	33.61	25	J**	All samples
	Vinyl Chloride	Acceptable	21.97	20	J**	All samples
	1,1-Dichloroethene	Acceptable	21.3	20	J**	All samples
	1,1,1-Trichloroethane	Acceptable	20.64	20	J	All samples
	Carbon Tetrachloride	Acceptable	26.3	20	J**	All samples
	Tetrachloroethene	Acceptable	20.48	20	J**	All samples
	2-Hexanone	Acceptable	48.57	40	J**	All samples

\*\*Qualification required for detected results only - associated results nondetect - no qualification required



ICV / CCV		RRF	%D	Limits	Qualifiers	Associated Samples	
ICV							
9/10/2020 2:17		Acceptable	Acceptable				
CCV							
9/10/2020 1:49		Vinyl Chloride	Acceptable	43.7	25	J / UJ	2009006-01, 2009006-07, 2009006-03RE1  ↓  2009006-01, 2009006-07, 2009006-03RE1
		Chloroethane	Acceptable	32.8	25	J / UJ	
		1,1-Dichloroethene	Acceptable	26.9	20	J / UJ	
		trans-1,2-Dichloroethene	Acceptable	25.7	20	J / UJ	
		1,1,1-Trichloroethane	Acceptable	43.5	25	J / UJ	
		Trichloroethene	Acceptable	25.1	20	J / UJ	
		Tetrachloroethene	Acceptable	30.7	20	J / UJ	
		Isopropylbenzene	Acceptable	25.1	25	J / UJ	
		1,2-Dibromo-3-chloropropane	Acceptable	58,5	30	J / UJ	
9/10/2020 8:38		Acceptable	Acceptable				
9/11/2020 9:55		Acceptable	Acceptable				
9/11/2020 7:06		Toluene	Acceptable	32.1	25	J / UJ	2009006-02RE1, 2009006-04RE1 through 2009006-06RE1, 2009007-06 through 2009007-08
9/15/2020 11:03		Dichlorodifluoromethane	Acceptable	40.1	40	J / UJ	2009007-06RE1, 2009007-09RE1 through 2009007-11RE1, 2009007-11RE2, 2009007-10RE2
9/15/2020 11:59			Acceptable	Acceptable			
9/16/2020 8:33		trans-1,3-Dichloropropene	Acceptable	21	20	J / UJ	2009007-08RE1, 2009007-09RE2

Tune  
Acceptable

MRL Check	%R	Limits	Qualifiers	Associated Samples
B20I010-MRL1	Acceptable			

Internal Standards	Area	Area Lower / Upper	Qualifiers	Associated Samples
		Limit		
		Acceptable		

#### Representativeness:

Were sampling procedures and design criteria met?

Were holding times met?

Was preservation criteria met? (0° C - 6° C)

Were Chain-of-Custody records complete and provided in data package?

Comments (note deviations): The cooler temperatures were 1.1 & 3.8 ° C.

Yes No N/A

Yes

Yes

Yes

Yes

Preservation	Cooler Temperature (Degrees C)	Preservation Criteria	Qualifier	Associated Samples
	Acceptable			

Holding Times	Analyte	Days to Extraction	HT Criteria	Qualifier	Associated Samples
		Acceptable			

#### Comparability:

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

Yes No N/A

Yes

#### Completeness (90%):

Are all data in this SDG usable?

Comments (note deviations):

Yes No N/A

Yes



**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):**Yes No N/A****Yes****Yes****Comment:**

As stated in the case narrative, all field samples were run at a 50x screening dilution and subsequent dilutions followed. Analytes are reported from the lowest sample dilution in which they were detected within the calibration range and reporting limits are raised accordingly.

As stated in the case narrative, as a result of the high concentrations of toluene, ethylbenzene, and m+p-xylene present in numerous field samples, carryover occurred in several instances in the project

Data is usable with appropriate qualifiers applied.

Data Validator:

Kristine MolloyDate: 1/21/2021

Data Reviewer:

Cherie ZakowskiDate: 1/25/2021





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB001-200909 (2009006-01)

Matrix: Water

Sampled: Sep-09-20 07:30

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B20I010	Sep-10-20	Sep-10-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"





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Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB001-200909 (2009006-01)

Matrix: Water

Sampled: Sep-09-20 07:30

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Chlorobenzene	U			2.00	ug/L	1	B201010	Sep-10-20	Sep-10-20
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.76			97.0%		73-124	"	"	"
1,2-Dichloroethane-d4	9.97			98.9%		84-122	"	"	"
Toluene-d8	9.68			96.8%		88-108	"	"	"
4-Bromofluorobenzene	9.66			96.6%		84-108	"	"	"





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Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004B-200909 (2009006-02RE1)

Matrix: Water

Sampled: Sep-09-20 16:30

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B20I012	Sep-11-20	Sep-11-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	5.34			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	4.93			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004B-200909 (2009006-02RE1)

Matrix: Water

Sampled: Sep-09-20 16:30

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.96			99.0%		73-124	"	"	"
1,2-Dichloroethane-d4	10.1			100%		84-122	"	"	"
Toluene-d8	9.85			98.5%		88-108	"	"	"
4-Bromofluorobenzene	9.74			97.4%		84-108	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW005-200909 (2009006-03RE1)

Matrix: Water

Sampled: Sep-09-20 13:40

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B201010	Sep-10-20	Sep-10-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	9.11			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	5.56			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW005-200909 (2009006-03RE1)

Matrix: Water

Sampled: Sep-09-20 13:40

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B201010	Sep-10-20	Sep-10-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.2			101%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			102%		84-122	"	"	"
Toluene-d8	9.66			96.6%		88-108	"	"	"
4-Bromofluorobenzene	9.72			97.2%		84-108	"	"	"





## Environmental Protection Agency Region 5

### US EPA Region 5 LSASD Analytical Services Branch

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Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

### Volatiles by GC/MS, EPA 8260C (modified) US EPA Region 5 LSASD Analytical Services Branch

A11-MW006-200909 (2009006-04RE1)

Matrix: Water

Sampled: Sep-09-20 11:45

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B20I012	Sep-11-20	Sep-11-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	2.28			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW006-200909 (2009006-04RE1)

Matrix: Water

Sampled: Sep-09-20 11:45

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.3			102%		73-124	"	"	"
1,2-Dichloroethane-d4	9.98			99.0%		84-122	"	"	"
Toluene-d8	9.77			97.7%		88-108	"	"	"
4-Bromofluorobenzene	9.66			96.6%		84-108	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW130A-200909 (2009006-05RE1)

Matrix: Water

Sampled: Sep-09-20 09:55

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B20I012	Sep-11-20	Sep-11-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	4.11			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	3.51			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW130A-200909 (2009006-05RE1)

Matrix: Water

Sampled: Sep-09-20 09:55

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.1			101%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			101%		84-122	"	"	"
Toluene-d8	9.79			97.9%		88-108	"	"	"
4-Bromofluorobenzene	8.94			89.4%		84-108	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW001-200909 (2009006-06RE1)

Matrix: Water

Sampled: Sep-09-20 16:25

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B20I012	Sep-11-20	Sep-11-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	5.16			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	7.58			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	2.41			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW001-200909 (2009006-06RE1)

Matrix: Water

Sampled: Sep-09-20 16:25

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.77			97.2%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			101%		84-122	"	"	"
Toluene-d8	9.65			96.5%		88-108	"	"	"
4-Bromofluorobenzene	9.70			97.0%		84-108	"	"	"





## Environmental Protection Agency Region 5

### US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Oct-23-20 13:16

### Volatiles by GC/MS, EPA 8260C (modified) US EPA Region 5 LSASD Analytical Services Branch

A11-FB001-200909 (2009006-07)

Matrix: Water

Sampled: Sep-09-20 18:00

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B201010	Sep-10-20	Sep-10-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-FB001-200909 (2009006-07)

Matrix: Water

Sampled: Sep-09-20 18:00

Received: Sep-10-20 10:05

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B201010	Sep-10-20	Sep-10-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.2			101%		73-124	"	"	"
1,2-Dichloroethane-d4	9.97			98.9%		84-122	"	"	"
Toluene-d8	9.90			99.0%		88-108	"	"	"
4-Bromofluorobenzene	9.50			95.0%		84-108	"	"	"





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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200910 (2009007-06)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethylbenzene	2630			100	ug/L	50	B20I012	Sep-11-20	Sep-11-20
m+p-Xylene	7600			200	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.60			95.4%		73-124	"	"	"
1,2-Dichloroethane-d4	10.4			103%		84-122	"	"	"
Toluene-d8	9.50			95.0%		88-108	"	"	"
4-Bromofluorobenzene	9.92			99.2%		84-108	"	"	"

A11-MW007-200910 (2009007-06RE1)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), (LCS), J		10.0	ug/L	5	B20I014	Sep-15-20	Sep-15-20
Chloromethane	U	(ICAL), J		10.0	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		10.0	"	"	"	"	"
Bromomethane	U			10.0	"	"	"	"	"
Chloroethane	U	(ICAL), J		10.0	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		10.0	"	"	"	"	"
1,1-Dichloroethene	U			10.0	"	"	"	"	"
Acetone	U			62.5	"	"	"	"	"
Carbon disulfide	U			10.0	"	"	"	"	"
Methylene chloride	U			10.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			10.0	"	"	"	"	"
1,1-Dichloroethane	U			10.0	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		10.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			10.0	"	"	"	"	"
2-Butanone	U			62.5	"	"	"	"	"
Bromochloromethane	U			10.0	"	"	"	"	"
Chloroform	U			10.0	"	"	"	"	"
1,1,1-Trichloroethane	U			10.0	"	"	"	"	"
Carbon tetrachloride	U			10.0	"	"	"	"	"
1,1-Dichloropropene	U			10.0	"	"	"	"	"
Benzene	U			10.0	"	"	"	"	"
1,2-Dichloroethane	U			10.0	"	"	"	"	"
Trichloroethene	U			10.0	"	"	"	"	"
1,2-Dichloropropane	U			10.0	"	"	"	"	"





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77 West Jackson Boulevard  
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Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200910 (2009007-06RE1)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dibromomethane	U			10.0	ug/L	5	B20I014	Sep-15-20	Sep-15-20
Bromodichloromethane	U			10.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			10.0	"	"	"	"	"
4-Methyl-2-pentanone	U			25.0	"	"	"	"	"
Toluene	U			10.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			10.0	"	"	"	"	"
1,1,2-Trichloroethane	U			10.0	"	"	"	"	"
Tetrachloroethene	U			10.0	"	"	"	"	"
1,3-Dichloropropane	U			10.0	"	"	"	"	"
2-Hexanone	U			25.0	"	"	"	"	"
Dibromochloromethane	U			10.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			10.0	"	"	"	"	"
Chlorobenzene	U			10.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			10.0	"	"	"	"	"
o-Xylene	U			10.0	"	"	"	"	"
Styrene	U			10.0	"	"	"	"	"
Bromoform	U			10.0	"	"	"	"	"
Isopropylbenzene	86.1			10.0	"	"	"	"	"
Bromobenzene	U			10.0	"	"	"	"	"
1,2,3-Trichloropropane	U			10.0	"	"	"	"	"
n-Propylbenzene	82.4			10.0	"	"	"	"	"
2-Chlorotoluene	U			10.0	"	"	"	"	"
1,3,5-Trimethylbenzene	11.1			10.0	"	"	"	"	"
4-Chlorotoluene	U			10.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			10.0	"	"	"	"	"
tert-Butylbenzene	U			10.0	"	"	"	"	"
1,2,4-Trimethylbenzene	53.5			10.0	"	"	"	"	"
sec-Butylbenzene	10.8			10.0	"	"	"	"	"
1,3-Dichlorobenzene	U			10.0	"	"	"	"	"
p-Isopropyltoluene	U			10.0	"	"	"	"	"
1,4-Dichlorobenzene	U			10.0	"	"	"	"	"
1,2-Dichlorobenzene	U			10.0	"	"	"	"	"
n-Butylbenzene	11.3			10.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			10.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			10.0	"	"	"	"	"
Hexachlorobutadiene	U			10.0	"	"	"	"	"
Naphthalene	11.4			10.0	"	"	"	"	"





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77 West Jackson Boulevard  
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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200910 (2009007-06RE1)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2,3-Trichlorobenzene	U			10.0	ug/L	5	B201014	Sep-15-20	Sep-15-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.74			96.8%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			101%		84-122	"	"	"
Toluene-d8	9.50			95.0%		88-108	"	"	"
4-Bromofluorobenzene	9.94			99.4%		84-108	"	"	"

A11-TB002-200910 (2009007-07)

Matrix: Water

Sampled: Sep-10-20 08:00

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), J		2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
Chloromethane	U	(ICAL), J		2.00	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U	(ICAL), J		2.00	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"





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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB002-200910 (2009007-07)

Matrix: Water

Sampled: Sep-10-20 08:00

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
cis-1,3-Dichloropropene	U			2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"





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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB002-200910 (2009007-07)

Matrix: Water

Sampled: Sep-10-20 08:00

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2,3-Trichlorobenzene	U			2.00	ug/L	1	B201012	Sep-11-20	Sep-11-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.70			96.4%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			101%		84-122	"	"	"
Toluene-d8	9.62			96.2%		88-108	"	"	"
4-Bromofluorobenzene	9.55			95.5%		84-108	"	"	"

A11-MW007-200910-D (2009007-08)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethylbenzene	2680			100	ug/L	50	B201012	Sep-11-20	Sep-11-20
m+p-Xylene	7920			200	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.92			98.6%		73-124	"	"	"
1,2-Dichloroethane-d4	9.93			98.5%		84-122	"	"	"
Toluene-d8	9.75			97.5%		88-108	"	"	"
4-Bromofluorobenzene	9.79			97.9%		84-108	"	"	"

A11-MW007-200910-D (2009007-08RE1)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), (LCS), J		10.0	ug/L	5	B201014	Sep-15-20	Sep-16-20
Chloromethane	U	(ICAL), J		10.0	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		10.0	"	"	"	"	"
Bromomethane	U			10.0	"	"	"	"	"
Chloroethane	U	(ICAL), J		10.0	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		10.0	"	"	"	"	"
1,1-Dichloroethene	U			10.0	"	"	"	"	"
Acetone	U			62.5	"	"	"	"	"
Carbon disulfide	U			10.0	"	"	"	"	"
Methylene chloride	U			10.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			10.0	"	"	"	"	"
1,1-Dichloroethane	U			10.0	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		10.0	"	"	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200910-D (2009007-08RE1)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
cis-1,2-Dichloroethene	U			10.0	ug/L	5	B201014	Sep-15-20	Sep-16-20
2-Butanone	U			62.5	"	"	"	"	"
Bromochloromethane	U			10.0	"	"	"	"	"
Chloroform	U			10.0	"	"	"	"	"
1,1,1-Trichloroethane	U			10.0	"	"	"	"	"
Carbon tetrachloride	U			10.0	"	"	"	"	"
1,1-Dichloropropene	U			10.0	"	"	"	"	"
Benzene	U			10.0	"	"	"	"	"
1,2-Dichloroethane	U			10.0	"	"	"	"	"
Trichloroethene	U			10.0	"	"	"	"	"
1,2-Dichloropropane	U			10.0	"	"	"	"	"
Dibromomethane	U			10.0	"	"	"	"	"
Bromodichloromethane	U			10.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			10.0	"	"	"	"	"
4-Methyl-2-pentanone	U			25.0	"	"	"	"	"
Toluene	U			10.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			10.0	"	"	"	"	"
1,1,2-Trichloroethane	U			10.0	"	"	"	"	"
Tetrachloroethene	U			10.0	"	"	"	"	"
1,3-Dichloropropane	U			10.0	"	"	"	"	"
2-Hexanone	U			25.0	"	"	"	"	"
Dibromochloromethane	U			10.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			10.0	"	"	"	"	"
Chlorobenzene	U			10.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			10.0	"	"	"	"	"
o-Xylene	U			10.0	"	"	"	"	"
Styrene	U			10.0	"	"	"	"	"
Bromoform	U			10.0	"	"	"	"	"
Isopropylbenzene	89.1			10.0	"	"	"	"	"
Bromobenzene	U			10.0	"	"	"	"	"
1,2,3-Trichloropropane	U			10.0	"	"	"	"	"
n-Propylbenzene	84.7			10.0	"	"	"	"	"
2-Chlorotoluene	U			10.0	"	"	"	"	"
1,3,5-Trimethylbenzene	11.6			10.0	"	"	"	"	"
4-Chlorotoluene	U			10.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			10.0	"	"	"	"	"
tert-Butylbenzene	U			10.0	"	"	"	"	"





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77 West Jackson Boulevard  
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Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW007-200910-D (2009007-08RE1)

Matrix: Water

Sampled: Sep-10-20 10:55

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2,4-Trimethylbenzene	55.7			10.0	ug/L	5	B201014	Sep-15-20	Sep-16-20
sec-Butylbenzene	11.3			10.0	"	"	"	"	"
1,3-Dichlorobenzene	U			10.0	"	"	"	"	"
p-Isopropyltoluene	U			10.0	"	"	"	"	"
1,4-Dichlorobenzene	U			10.0	"	"	"	"	"
1,2-Dichlorobenzene	U			10.0	"	"	"	"	"
n-Butylbenzene	12.4			10.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			10.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			10.0	"	"	"	"	"
Hexachlorobutadiene	U			10.0	"	"	"	"	"
Naphthalene	13.2			10.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			10.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.84			97.8%		73-124	"	"	"
1,2-Dichloroethane-d4	10.3			102%		84-122	"	"	"
Toluene-d8	9.49			94.9%		88-108	"	"	"
4-Bromofluorobenzene	10.0			100%		84-108	"	"	"

A11-MW002-200910 (2009007-09RE1)

Matrix: Water

Sampled: Sep-10-20 13:05

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Toluene	39300			1600	ug/L	800	B201014	Sep-15-20	Sep-15-20
Ethylbenzene	8260			1600	"	"	"	"	"
m+p-Xylene	26000			3200	"	"	"	"	"
o-Xylene	6820			1600	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.62			95.6%		73-124	"	"	"
1,2-Dichloroethane-d4	9.94			98.6%		84-122	"	"	"
Toluene-d8	9.49			94.9%		88-108	"	"	"
4-Bromofluorobenzene	9.48			94.8%		84-108	"	"	"





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77 West Jackson Boulevard  
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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW002-200910 (2009007-09RE2)

Matrix: Water

Sampled: Sep-10-20 13:05

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), (LCS), J		50.0	ug/L	25	B201014	Sep-15-20	Sep-16-20
Chloromethane	U	(ICAL), J		50.0	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		50.0	"	"	"	"	"
Bromomethane	U			50.0	"	"	"	"	"
Chloroethane	U	(ICAL), J		50.0	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		50.0	"	"	"	"	"
1,1-Dichloroethene	U			50.0	"	"	"	"	"
Acetone	U			312	"	"	"	"	"
Carbon disulfide	U			50.0	"	"	"	"	"
Methylene chloride	U			50.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			50.0	"	"	"	"	"
1,1-Dichloroethane	U			50.0	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		50.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			50.0	"	"	"	"	"
2-Butanone	U			312	"	"	"	"	"
Bromochloromethane	U			50.0	"	"	"	"	"
Chloroform	U			50.0	"	"	"	"	"
1,1,1-Trichloroethane	U			50.0	"	"	"	"	"
Carbon tetrachloride	U			50.0	"	"	"	"	"
1,1-Dichloropropene	U			50.0	"	"	"	"	"
Benzene	U			50.0	"	"	"	"	"
1,2-Dichloroethane	U			50.0	"	"	"	"	"
Trichloroethene	U			50.0	"	"	"	"	"
1,2-Dichloropropane	U			50.0	"	"	"	"	"
Dibromomethane	U			50.0	"	"	"	"	"
Bromodichloromethane	U			50.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			50.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			50.0	"	"	"	"	"
1,1,2-Trichloroethane	U			50.0	"	"	"	"	"
Tetrachloroethene	U			50.0	"	"	"	"	"
1,3-Dichloropropane	U			50.0	"	"	"	"	"
2-Hexanone	U			125	"	"	"	"	"
Dibromochloromethane	U			50.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			50.0	"	"	"	"	"
Chlorobenzene	U			50.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			50.0	"	"	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

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Project Manager: Terese Van Donsel

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Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW002-200910 (2009007-09RE2)

Matrix: Water

Sampled: Sep-10-20 13:05

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Styrene	U			50.0	ug/L	25	B201014	Sep-15-20	Sep-16-20
Bromoform	U			50.0	"	"	"	"	"
Isopropylbenzene	90.0			50.0	"	"	"	"	"
Bromobenzene	U			50.0	"	"	"	"	"
1,2,3-Trichloropropane	U			50.0	"	"	"	"	"
n-Propylbenzene	129			50.0	"	"	"	"	"
2-Chlorotoluene	U			50.0	"	"	"	"	"
1,3,5-Trimethylbenzene	202			50.0	"	"	"	"	"
4-Chlorotoluene	U			50.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			50.0	"	"	"	"	"
tert-Butylbenzene	U			50.0	"	"	"	"	"
1,2,4-Trimethylbenzene	622			50.0	"	"	"	"	"
sec-Butylbenzene	U			50.0	"	"	"	"	"
1,3-Dichlorobenzene	U			50.0	"	"	"	"	"
p-Isopropyltoluene	U			50.0	"	"	"	"	"
1,4-Dichlorobenzene	U			50.0	"	"	"	"	"
1,2-Dichlorobenzene	U			50.0	"	"	"	"	"
n-Butylbenzene	U			50.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			50.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			50.0	"	"	"	"	"
Hexachlorobutadiene	U			50.0	"	"	"	"	"
Naphthalene	55.2			50.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			50.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.36			93.1%		73-124	"	"	"
1,2-Dichloroethane-d4	10.1			100%		84-122	"	"	"
Toluene-d8	9.73			97.3%		88-108	"	"	"
4-Bromofluorobenzene	9.71			97.1%		84-108	"	"	"





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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004A-200910 (2009007-10RE1)

Matrix: Water

Sampled: Sep-10-20 15:50

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Toluene</b>	<b>42600</b>			1000	ug/L	500	B20I014	Sep-15-20	Sep-15-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>Dibromofluoromethane</i>	9.68			96.3%		73-124	"	"	"
<i>1,2-Dichloroethane-d4</i>	10.4			103%		84-122	"	"	"
<i>Toluene-d8</i>	9.69			96.9%		88-108	"	"	"
<i>4-Bromofluorobenzene</i>	8.91			89.1%		84-108	"	"	"

A11-MW004A-200910 (2009007-10RE2)

Matrix: Water

Sampled: Sep-10-20 15:50

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Dichlorodifluoromethane</b>	U	(ICAL), (LCS), J		50.0	ug/L	25	B20I014	Sep-15-20	Sep-15-20
<b>Chloromethane</b>	U	(ICAL), J		50.0	"	"	"	"	"
<b>Vinyl chloride</b>	U	(ICAL), J		50.0	"	"	"	"	"
<b>Bromomethane</b>	U			50.0	"	"	"	"	"
<b>Chloroethane</b>	U	(ICAL), J		50.0	"	"	"	"	"
<b>Trichlorofluoromethane</b>	U	(ICAL), J		50.0	"	"	"	"	"
<b>1,1-Dichloroethene</b>	U			50.0	"	"	"	"	"
<b>Acetone</b>	U			312	"	"	"	"	"
<b>Carbon disulfide</b>	U			50.0	"	"	"	"	"
<b>Methylene chloride</b>	U			50.0	"	"	"	"	"
<b>trans-1,2-Dichloroethene</b>	U			50.0	"	"	"	"	"
<b>1,1-Dichloroethane</b>	U			50.0	"	"	"	"	"
<b>2,2-Dichloropropane</b>	U	(LCS), J		50.0	"	"	"	"	"
<b>cis-1,2-Dichloroethene</b>	U			50.0	"	"	"	"	"
<b>2-Butanone</b>	U			312	"	"	"	"	"
<b>Bromochloromethane</b>	U			50.0	"	"	"	"	"
<b>Chloroform</b>	U			50.0	"	"	"	"	"
<b>1,1,1-Trichloroethane</b>	U			50.0	"	"	"	"	"
<b>Carbon tetrachloride</b>	U			50.0	"	"	"	"	"
<b>1,1-Dichloropropene</b>	U			50.0	"	"	"	"	"
<b>Benzene</b>	U			50.0	"	"	"	"	"
<b>1,2-Dichloroethane</b>	U			50.0	"	"	"	"	"
<b>Trichloroethene</b>	U			50.0	"	"	"	"	"
<b>1,2-Dichloropropane</b>	U			50.0	"	"	"	"	"
<b>Dibromomethane</b>	U			50.0	"	"	"	"	"





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77 West Jackson Boulevard  
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Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004A-200910 (2009007-10RE2)

Matrix: Water

Sampled: Sep-10-20 15:50

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Bromodichloromethane	U			50.0	ug/L	25	B201014	Sep-15-20	Sep-15-20
cis-1,3-Dichloropropene	U			50.0	"	"	"	"	"
4-Methyl-2-pentanone	U			125	"	"	"	"	"
trans-1,3-Dichloropropene	U			50.0	"	"	"	"	"
1,1,2-Trichloroethane	U			50.0	"	"	"	"	"
Tetrachloroethene	U			50.0	"	"	"	"	"
1,3-Dichloropropane	U			50.0	"	"	"	"	"
2-Hexanone	U			125	"	"	"	"	"
Dibromochloromethane	U			50.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			50.0	"	"	"	"	"
Chlorobenzene	U			50.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			50.0	"	"	"	"	"
Ethylbenzene	365			50.0	"	"	"	"	"
m+p-Xylene	538			100	"	"	"	"	"
o-Xylene	66.6			50.0	"	"	"	"	"
Styrene	U			50.0	"	"	"	"	"
Bromoform	U			50.0	"	"	"	"	"
Isopropylbenzene	U			50.0	"	"	"	"	"
Bromobenzene	U			50.0	"	"	"	"	"
1,2,3-Trichloropropane	U			50.0	"	"	"	"	"
n-Propylbenzene	U			50.0	"	"	"	"	"
2-Chlorotoluene	U			50.0	"	"	"	"	"
1,3,5-Trimethylbenzene	U			50.0	"	"	"	"	"
4-Chlorotoluene	U			50.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			50.0	"	"	"	"	"
tert-Butylbenzene	U			50.0	"	"	"	"	"
1,2,4-Trimethylbenzene	U			50.0	"	"	"	"	"
sec-Butylbenzene	U			50.0	"	"	"	"	"
1,3-Dichlorobenzene	U			50.0	"	"	"	"	"
p-Isopropyltoluene	U			50.0	"	"	"	"	"
1,4-Dichlorobenzene	U			50.0	"	"	"	"	"
1,2-Dichlorobenzene	U			50.0	"	"	"	"	"
n-Butylbenzene	U			50.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			50.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			50.0	"	"	"	"	"
Hexachlorobutadiene	U			50.0	"	"	"	"	"
Naphthalene	U			50.0	"	"	"	"	"





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Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004A-200910 (2009007-10RE2)

Matrix: Water

Sampled: Sep-10-20 15:50

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2,3-Trichlorobenzene	U			50.0	ug/L	25	B20I014	Sep-15-20	Sep-15-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.35			92.9%		73-124	"	"	"
1,2-Dichloroethane-d4	10.1			101%		84-122	"	"	"
Toluene-d8	9.49			94.9%		88-108	"	"	"
4-Bromofluorobenzene	9.68			96.8%		84-108	"	"	"

A11-MW003-200910 (2009007-11RE1)

Matrix: Water

Sampled: Sep-10-20 08:45

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
m+p-Xylene	2430			100	ug/L	25	B20I014	Sep-15-20	Sep-15-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.95			98.9%		73-124	"	"	"
1,2-Dichloroethane-d4	10.4			103%		84-122	"	"	"
Toluene-d8	9.44			94.4%		88-108	"	"	"
4-Bromofluorobenzene	9.63			96.3%		84-108	"	"	"

A11-MW003-200910 (2009007-11RE2)

Matrix: Water

Sampled: Sep-10-20 08:45

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U	(ICAL), (LCS), J		10.0	ug/L	5	B20I014	Sep-15-20	Sep-15-20
Chloromethane	U	(ICAL), J		10.0	"	"	"	"	"
Vinyl chloride	U	(ICAL), J		10.0	"	"	"	"	"
Bromomethane	U			10.0	"	"	"	"	"
Chloroethane	U	(ICAL), J		10.0	"	"	"	"	"
Trichlorofluoromethane	U	(ICAL), J		10.0	"	"	"	"	"
1,1-Dichloroethene	U			10.0	"	"	"	"	"
Acetone	U			62.5	"	"	"	"	"
Carbon disulfide	U			10.0	"	"	"	"	"
Methylene chloride	U			10.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			10.0	"	"	"	"	"
1,1-Dichloroethane	U			10.0	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		10.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			10.0	"	"	"	"	"





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Oct-23-20 13:16

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-MW003-200910 (2009007-11RE2)

Matrix: Water

Sampled: Sep-10-20 08:45

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
2-Butanone	U			62.5	ug/L	5	B20I014	Sep-15-20	Sep-15-20
Bromochloromethane	U			10.0	"	"	"	"	"
Chloroform	U			10.0	"	"	"	"	"
1,1,1-Trichloroethane	U			10.0	"	"	"	"	"
Carbon tetrachloride	U			10.0	"	"	"	"	"
1,1-Dichloropropene	U			10.0	"	"	"	"	"
Benzene	U			10.0	"	"	"	"	"
1,2-Dichloroethane	U			10.0	"	"	"	"	"
Trichloroethene	U			10.0	"	"	"	"	"
1,2-Dichloropropane	U			10.0	"	"	"	"	"
Dibromomethane	U			10.0	"	"	"	"	"
Bromodichloromethane	U			10.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			10.0	"	"	"	"	"
4-Methyl-2-pentanone	U			25.0	"	"	"	"	"
Toluene	U			10.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			10.0	"	"	"	"	"
1,1,2-Trichloroethane	U			10.0	"	"	"	"	"
Tetrachloroethene	U			10.0	"	"	"	"	"
1,3-Dichloropropane	U			10.0	"	"	"	"	"
2-Hexanone	U			25.0	"	"	"	"	"
Dibromochloromethane	U			10.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			10.0	"	"	"	"	"
Chlorobenzene	U			10.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			10.0	"	"	"	"	"
Ethylbenzene	201			10.0	"	"	"	"	"
o-Xylene	U			10.0	"	"	"	"	"
Styrene	U			10.0	"	"	"	"	"
Bromoform	U			10.0	"	"	"	"	"
Isopropylbenzene	28.3			10.0	"	"	"	"	"
Bromobenzene	U			10.0	"	"	"	"	"
1,2,3-Trichloropropane	U			10.0	"	"	"	"	"
n-Propylbenzene	28.2			10.0	"	"	"	"	"
2-Chlorotoluene	U			10.0	"	"	"	"	"
1,3,5-Trimethylbenzene	34.8			10.0	"	"	"	"	"
4-Chlorotoluene	U			10.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			10.0	"	"	"	"	"
tert-Butylbenzene	U			10.0	"	"	"	"	"





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Project: SE Rockford GW Contamination  
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Reported:  
Oct-23-20 13:16

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW003-200910 (2009007-11RE2)

Matrix: Water

Sampled: Sep-10-20 08:45

Received: Sep-11-20 10:10

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2,4-Trimethylbenzene	113			10.0	ug/L	5	B201014	Sep-15-20	Sep-15-20
sec-Butylbenzene	12.0			10.0	"	"	"	"	"
1,3-Dichlorobenzene	U			10.0	"	"	"	"	"
p-Isopropyltoluene	U			10.0	"	"	"	"	"
1,4-Dichlorobenzene	U			10.0	"	"	"	"	"
1,2-Dichlorobenzene	U			10.0	"	"	"	"	"
n-Butylbenzene	U			10.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			10.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			10.0	"	"	"	"	"
Hexachlorobutadiene	U			10.0	"	"	"	"	"
Naphthalene	U			10.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			10.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.82			97.6%		73-124	"	"	"
1,2-Dichloroethane-d4	10.4			103%		84-122	"	"	"
Toluene-d8	9.57			95.7%		88-108	"	"	"
4-Bromofluorobenzene	10.1			101%		84-108	"	"	"



## **December 2020 Data Validation Reports and Data Packages**



**Techlaw Document Controlled Number: 83139-1-23-612-DV-0016**  
UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V  
SUPERFUND AND EMERGENCY MANAGEMENT DIVISION

DATE:

SUBJECT: Review of Data  
Received for Review on: January 11, 2021

FROM: Allison C Harvet, TechLaw Consultants, Inc.  
Contractor, Environmental Services Assistance Team (ESAT)

THROUGH: Michelle Kerr  
Region 5 ESAT Contracting Officer's Representative

TO: Data User: CDM Smith  
Contact Person: John Grabs  
Email address: grabsjc@cdmsmith.com

**Stage\_2B\_Validation\_Electronic\_And\_Manual (S2BVEM) Data Review Narrative**

We have reviewed the data for the following case:

SITE Name: Southeast Rockford Groundwater, Area 11 (IL)

Case No: 49238 MA No: \_\_\_\_\_ SDG No: E3YH7

Number and Type of Samples: 11 waters (SVOA SIM)

Sample Numbers: E3YH1 – E3YH5, E3YH7 – E3YH9, E3YJ0 – E3YJ2

Laboratory: Pace Analytical Services, LLC Hrs. for Review:

Following are our findings:



**Below is a summary of the out-of-control audits and the possible effects on the data for this case:**

Eleven (11) preserved water samples labeled E3YH1 through E3YH5, E3YH7 through E3YH9, and E3YJ0 through E3YJ2, were shipped to Pace Analytical Services LLC (EQI) located in West Columbia, SC. The samples were collected 12/01-02/2020 and received 12/02/2020 and 12/03/2020 intact. Four (4) samples; E3YH7, E3YJ0, E3YJ1 and E3YJ2, were received at the elevated temperature of 6.1°C. The remaining samples arrived properly cooled between 2.6°C and 4.1°C. All samples were analyzed according to CLP SOW SOM02.4, [Oct 2016] (and MA: 3054.0 – 1,4-Dioxane Analysis with Lower CRQL) and reviewed according to the QAPP, the September 2017 NFG for SOM02.4 (EPA-540-R-2017-002) and the Region 5 Organic CLP Validation SOP, DCN/SOP 83074-8-33-601-SO-1143.

Sample E3YH9 was designated by the samplers to be used for laboratory QC, i.e. MS/MSD analyses.

Sample E3YJ2 was identified as a field blank. Sample E3YH1 was identified as a field duplicate of sample E3YH2.

The sample results have been reviewed for compliance with the QAPP worksheets and all non-compliance are described in Section 17. – QAPP Compliance

‘Only outliers and non-compliances are discussed in the narrative’.

**1. PRESERVATION AND HOLDING TIMES**

NONE FOUND.

**2. GC/MS and GC/ECD INSTRUMENT PERFORMANCE CHECK**

NONE FOUND.

**3. INITIAL CALIBRATION**

NONE FOUND.

**4. INITIAL CALIBRATION VERIFICATION**

NONE FOUND.

**5. CONTINUING CALIBRATION**

NONE FOUND.

**6. BLANKS**

NONE FOUND.

**7. DEUTERATED MONITORING COMPOUNDS / SURROGATES**

NONE FOUND.

**8. MATRIX SPIKE/MATRIX SPIKE DUPLICATE**

The following samples reported percent recovery below the QC criteria specified in MA: 3054.0 (15-120 %R). Detects in the unspiked sample, E3HY9(DL) is qualified as estimated J.

E3YH9MSD  
1,4-Dioxane

The relative percent difference (RPD) between the following samples is outside the QC criteria specified in MA: 3054.0 (0-50 RPD). Detects in the unspiked sample, E3YH9(DL) is qualified as estimated J.

E3YH9MS, E3YH9MSD  
1,4-Dioxane

**9. CLEANUP PROCEDURES**



NONE FOUND.

**10. LABORATORY CONTROL SAMPLE**

NONE FOUND.

**11. INTERNAL STANDARD**

NONE FOUND.

**12. TARGET ANALYTE QUANTITATION LIMIT**

**Method – Semivolatiles by SIM**

EXES-790

The following samples have analyte results greater than or equal to detection limit (MDL) and below quantitation limit (CRQL). Detects are qualified as estimated J.

E3YH2, E3YH3  
1,4-Dioxane

**13. TENTATIVELY IDENTIFIED COMPOUNDS**

Not Validated for this Stage of Review.

**14. SYSTEM PERFORMANCE**

NONE FOUND.

**15. FIELD QC SAMPLES**

Review not required under specified validation stage.

**16. SAMPLE RESULTS**

The following samples reported analyte concentrations above the calibration range. No dilutions were performed as these samples are QC samples. Detects are qualified as estimated J.

E3YH9MS, E3YH9MSD  
1,4-Dioxane

**17. QAPP COMPLIANCE**

The analytical package fulfilled the QAPP QC components requirements identified in the Southeast Rockford GW QAPP – Area 11.

Case No: 49238

Site Name: Southeast Rockford Groundwater, Area 11 (IL)

Page 5 of 6

SDG No: E3YH5

Laboratory: Pace (EQI)

The raw data package was missing the Form 3 for QC sample SLCS76.



## Validation Data Qualifier Sheet

QualifiersData Qualifier Definitions

U	The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
J	The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
J+	The result is an estimated quantity, but the results may be biased high.
J-	The result is an estimated quantity, but the results may be biased low.
NJ	The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
UJ	The analyte was analyzed for, but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
R	The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.
C	The Target Pesticide or Aroclor analyte identification has been confirmed by Gas Chromatograph/Mass Spectrometer (GC/MS).
X	The Target Pesticide or Aroclor analyte identification was not confirmed when GC/MS analysis was performed.

# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD GROUND WATER CONTAMINATION Project	GroupID: 49238/EPW14035/E3YH7	Lab Name: Pace Analytical Services, LLC
--	-------------------------------	---

Sample Number: E3YH1	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW007	pH: 8	Sample Date: 12/02/2020	Sample Time: 10:50:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	0.19	U	ug/L	0.19	U	1.0	YES	S2BVEM



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH2	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW007	pH: 8	Sample Date: 12/02/2020	Sample Time: 10:50:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	0.069	J	ug/L	0.069	J	1.0	YES	S2BVEM

# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH3	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW004A	pH: 8	Sample Date: 12/02/2020	Sample Time: 15:00:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	0.15	J	ug/L	0.15	J	1.0	YES	S2BVEM



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH4	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW003	pH: 8	Sample Date: 12/02/2020	Sample Time: 08:45:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	4.6		ug/L	4.6	D	2.0	YES	S2BVEM

# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH5	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW002	pH: 8	Sample Date: 12/02/2020	Sample Time: 13:10:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	1.1		ug/L	1.1	D	2.0	YES	S2BVEM



# Sample Summary Report

**Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project**

**GroupID: 49238/EPW14035/E3YH7**

**Lab Name: Pace Analytical Services, LLC**

Sample Number: E3YH7	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW130A	pH: 8	Sample Date: 12/01/2020	Sample Time: 09:35:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	4.0		ug/L	4.0	D	2.0	YES	S2BVEM

# Sample Summary Report

**Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project**

**GroupID: 49238/EPW14035/E3YH7**

**Lab Name: Pace Analytical Services, LLC**

Sample Number: E3YH8	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW006	pH: 8	Sample Date: 12/01/2020	Sample Time: 12:11:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	4.1		ug/L	4.1	D	2.0	YES	S2BVEM



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH9	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW005	pH: 8	Sample Date: 12/01/2020	Sample Time: 13:50:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	4.5	J	ug/L	4.5	D	2.0	YES	S2BVEM

# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH9MS	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location:	pH: 8	Sample Date: 12/01/2020	Sample Time: 13:50:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Spike	5.6	J	ug/L	5.6	E	1.0	YES	S2BVEM



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YH9MSD	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location:	pH: 8	Sample Date: 12/01/2020	Sample Time: 13:50:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Spike	4.5	J	ug/L	4.5	E	1.0	YES	S2BVEM

# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YJ0	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW004B	pH: 8	Sample Date: 12/01/2020	Sample Time: 16:45:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	6.3		ug/L	6.3	D	5.0	YES	S2BVEM



# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YJ1	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-MW001	pH: 8	Sample Date: 12/01/2020	Sample Time: 15:31:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	5.0		ug/L	5.0	D	2.0	YES	S2BVEM

# Sample Summary Report

Project Name: SOUTHEAST ROCKFORD  
GROUND WATER CONTAMINATION Project

GroupID: 49238/EPW14035/E3YH7

Lab Name: Pace Analytical Services, LLC

Sample Number: E3YJ2	Method: Semivolatiles by SIM	Matrix: Water	MA Number: 3054.0
Sample Location: A11-FB001	pH: 8	Sample Date: 12/01/2020	Sample Time: 17:05:00
% Moisture:		% Solids: 0.0	

Analyte Name	Analyte Type	Validation Result	Validation Flag	Units	Lab Result	Lab Flag	Dilution Factor	Reportable	Validation Level
1,4-Dioxane	Target	0.19	U	ug/L	0.19	U	1.0	YES	S2BVEM



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 2012003

**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:** Groundwater

**Collection date:** 12/01/20

**Analysis/Methods:** Wet Chemistry: Anions - EPA 300.0

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>
2012003-01	A11-MW130A-201201
2012003-02	A11-MW006-201201
2012003-03	A11-MW005-201201
2012003-04	A11-MW004B-201201
2012003-05	A11-MW001-201201
2012003-06	A11-FB001-201201

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters (Anions 300.0)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?			N/A
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)			N/A
Laboratory Control Spike Duplicates RPD within limits?			N/A
Laboratory Duplicate RPDs within limits?			Yes
<u>Comments (note deviations):</u>			

<u>Field Duplicates</u>	<u>Sample</u>	<u>Duplicate</u>	<u>%RPD</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A					

---

<u>MS/MSD</u>	<u>%RPD</u>	<u>Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				

---

<u>LCS/LCSD</u>	<u>%RPD</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
N/A				

---

<u>Laboratory Duplicate</u>	<u>%RPD</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
E20L003-DUP1	Acceptable			

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)			Yes
Laboratory Control Sample criteria met?			Yes
Were the Laboratory Method Blank results all < RL?			Yes
Were the Field Blanks results all < RL?			Yes
Was the ICAL criteria met?			Yes
Was the CCV criteria met?			Yes
Was the Tuning criteria met?			N/A
Were the Surrogate % recoveries within laboratory determined control limits?			N/A
Were the Internal Standard areas within ± 50 - 150%?			N/A
<u>Comments (note deviations):</u>			

<u>Blanks</u>	<u>Concentration</u>	<u>MDL /PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
E20L003-BLK1				
Nitrogen, Nitrate	Nondetect			
Sulfate	Nondetect			

<b>ICB/CCB</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
ICB	Nitrogen, Nitrate Sulfate	Nondetect 0.03	0.10 / 0.12					
CCB1	Nitrogen, Nitrate Sulfate	Nondetect 0.03	0.10 / 0.12	None	Sample results nondetect or > RL			
<b>Field Blank</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
A11-FB001-201201		Nondetect						
<b>Surrogates</b>		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
N/A								
<b>MS/MSD</b>		<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
E20L003-MS1								
Nitrogen, Nitrate Sulfate		Acceptable 68%	80-120 80-120	J- / UJ	All samples			
<b>LCS/LCSD</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
E20L003-BS1								
Nitrogen, Nitrate Sulfate		Acceptable Acceptable	90-110 90-110					
<b>ICV</b>			<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
Nitrogen, Nitrate Sulfate			Acceptable Acceptable					
<b>CCV</b>			<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
Nitrogen, Nitrate Sulfate			Acceptable Acceptable					
<b>MRL Check</b>			<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
B20L003-MRL1								
Nitrogen, Nitrate Sulfate			Acceptable Acceptable					
<b>Tune</b>								
N/A								
<b>Internal Standards</b>		<b><u>Area</u></b>	<b><u>Area Lower / Upper Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
N/A								
<b>Representativeness:</b>						<b><u>Yes</u></b>	<b><u>No</u></b>	<b><u>N/A</u></b>
Were sampling procedures and design criteria met?							Yes	
Were holding times met?							Yes	
Was preservation criteria met? (0° C - 6° C)							Yes	
Were Chain-of-Custody records complete and provided in data package?							No	
<u>Comments (note deviations):</u> The cooler temperature was 0.3 ° C.								
<b>Preservation</b>		<b><u>Cooler Temperature (Degrees C)</u></b>	<b><u>Preservation Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>			
		Acceptable						
<b>Holding Times</b>		<b><u>Analyte</u></b>	<b><u>Days to Extraction</u></b>	<b><u>HT Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>		
			Acceptable					



**Comparability:**

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Completeness (90%):**

Are all data in this SDG usable?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Yes**

**Comment:**

Data is usable with appropriate qualifiers applied.

Data Validator:

*Kristine Molloy*

Date: 1/19/2021

Data Reviewer:

Cherie Zakowski

Date: 1/25/2021



**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Nov-09-20 15:25

**Anions by Ion Chromatography, EPA 300.0 (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW002-200910 (2009007-01)**

**Matrix: Water**

**Sampled: Sep-10-20 13:05**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	U			0.12	mg/L	1	B20I013	Sep-11-20	Sep-11-20
Nitrate - NO3	U			0.12	"	"	"	"	"

**A11-MW007-200910 (2009007-02)**

**Matrix: Water**

**Sampled: Sep-10-20 10:55**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	2.96			0.12	mg/L	1	B20I013	Sep-11-20	Sep-11-20
Nitrate - NO3	U			0.12	"	"	"	"	"

**A11-MW004A-200910 (2009007-03)**

**Matrix: Water**

**Sampled: Sep-10-20 15:50**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	33.4			0.12	mg/L	1	B20I013	Sep-11-20	Sep-11-20
Nitrate - NO3	1.93			0.12	"	"	"	"	"

**A11-MW007-200910-D (2009007-04)**

**Matrix: Water**

**Sampled: Sep-10-20 10:55**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	2.93			0.12	mg/L	1	B20I013	Sep-11-20	Sep-11-20
Nitrate - NO3	U			0.12	"	"	"	"	"

**A11-MW003-200910-D (2009007-05)**

**Matrix: Water**

**Sampled: Sep-10-20 08:45**

**Received: Sep-11-20 10:10**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	11.3			0.12	mg/L	1	B20I013	Sep-11-20	Sep-11-20
Nitrate - NO3	U			0.12	"	"	"	"	"





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**US EPA Region 5 LSASD Analytical Services Branch**

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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Nov-09-20 15:25

**Notes and Definitions**

U Not Detected  
NR Not Reported  
Q QC limit Exceeded

**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 2012005

**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:** Groundwater

**Collection date:** 12/02/20

**Analysis/Methods:** Wet Chemistry: Anions - EPA 300.0

**Samples in SDG:**

Lab ID	Sample Number
2012005-01	A11-MW007-201201-D
2012005-02	A11-MW007-201201
2012005-03	A11-MW004A-201201
2012005-04	A11-MW003-201201
2012005-05	A11-MW002-201201

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters (Anions 300.0)**

Precision:	Yes	No	N/A
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	N/A		
Laboratory Control Spike Duplicates RPD within limits?	N/A		
Laboratory Duplicate RPDs within limits?	Yes		
<u>Comments (note deviations):</u>			

Field Duplicates	Sample A11-MW007-201201	Duplicate A11-MW007-201201-D	%RPD	Qualifiers	Associated Samples
			Acceptable		

MS/MSD	%RPD	Limit	Qualifiers	Associated Samples
N/A				

LCS/LCSD	%RPD	Limits	Qualifiers	Associated Samples
N/A				

Laboratory Duplicate	%RPD	Limits	Qualifiers	Associated Samples
E20L006-DUP1	Acceptable			

Accuracy:	Yes	No	N/A
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	N/A		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

Blanks	Concentration	MDL /PQL	Qualifiers	Associated Samples
E20L006-BLK1				
Nitrogen, Nitrate	Nondetect			
Sulfate	Nondetect			



<b>ICB/CCB</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
ICB	Nitrogen, Nitrate Sulfate	Nondetect 0.03	0.1 / 0.12	None	Sample results > RL
CCB1	Nitrogen, Nitrate Sulfate	Nondetect 0.03	0.1 / 0.12	None	Sample results > RL
<b>Field Blank</b> N/A		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>Surrogates</b> N/A		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>MS/MSD</b> E20L006-MS1 Nitrogen, Nitrate Sulfate		<b><u>%R</u></b>  Acceptable Acceptable	<b><u>Limits (%)</u></b>  80-120 80-120	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>LCS/LCSD</b> E20L006-BS1 Nitrogen, Nitrate Sulfate		<b><u>%R</u></b>  Acceptable Acceptable	<b><u>Limits</u></b>  90-110 90-110	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>ICV</b>	Nitrogen, Nitrate Sulfate	<b><u>%R</u></b> Acceptable Acceptable	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>CCV</b>	Nitrogen, Nitrate Sulfate	<b><u>%R</u></b> Acceptable Acceptable	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>MRL Check</b> B20L006-MRL1 Nitrogen, Nitrate Sulfate		<b><u>%R</u></b>  Acceptable Acceptable	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>Tune</b> N/A					
<b>Internal Standards</b> N/A	<b><u>Area</u></b>	<b><u>Area Lower / Upper Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
<b>Representativeness:</b>					
Were sampling procedures and design criteria met?					<b>Yes</b>
Were holding times met?					<b>Yes</b>
Was preservation criteria met? (0° C - 6° C)					<b>Yes</b>
Were Chain-of-Custody records complete and provided in data package?					<b>No</b>
<u>Comments (note deviations):</u> The cooler temperature was 0.6 ° C.					
<b>Preservation</b>	<b><u>Cooler Temperature (Degrees C)</u></b> Acceptable	<b><u>Preservation Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>	
<b>Holding Times</b>	<b><u>Analyte</u></b>	<b><u>Days to Extraction</u></b> Acceptable	<b><u>HT Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>

**Comparability:**

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Completeness (90%):**

Are all data in this SDG usable?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Yes**

**Comment:**

Data is usable as reported.

Data Validator:

*Kristine Molloy*

Date: 1/20/2021

Data Reviewer:

Cherie Zakowski

Date: 1/23/2021





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Dec-07-20 16:18

**Anions by Ion Chromatography, EPA 300.0 (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW130A-201201 (2012003-01)**

**Matrix: Water**

**Sampled: Dec-01-20 09:35**

**Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	17.3			0.12	mg/L	1	B20L003	Dec-02-20	Dec-02-20
Nitrate - NO3	6.26			0.12	"	"	"	"	"

**A11-MW006-201201 (2012003-02)**

**Matrix: Water**

**Sampled: Dec-01-20 12:11**

**Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	5.56			0.12	mg/L	1	B20L003	Dec-02-20	Dec-02-20
Nitrate - NO3	U			0.12	"	"	"	"	"

**A11-MW005-201201 (2012003-03)**

**Matrix: Water**

**Sampled: Dec-01-20 13:50**

**Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	33.7	(MS), L		0.12	mg/L	1	B20L003	Dec-02-20	Dec-02-20
Nitrate - NO3	14.1			0.12	"	"	"	"	"

**A11-MW004B-201201 (2012003-04)**

**Matrix: Water**

**Sampled: Dec-01-20 16:45**

**Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	18.9			0.12	mg/L	1	B20L003	Dec-02-20	Dec-02-20
Nitrate - NO3	5.15			0.12	"	"	"	"	"

**A11-MW001-201201 (2012003-05)**

**Matrix: Water**

**Sampled: Dec-01-20 05:31**

**Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	30.4			0.12	mg/L	1	B20L003	Dec-02-20	Dec-02-20
Nitrate - NO3	11.4			0.12	"	"	"	"	"

**A11-FB001-201201 (2012003-06)**

**Matrix: Water**

**Sampled: Dec-01-20 17:01**

**Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO4	U			0.12	mg/L	1	B20L003	Dec-02-20	Dec-02-20
Nitrate - NO3	U			0.12	"	"	"	"	"



**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
Phone: (312) 353-8370 Fax: (312) 886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Dec-07-20 16:18

**Notes and Definitions**

- L The identification of the analyte is acceptable; the reported value may be biased low. The actual value is expected to be greater than the reported value.
- (MS) Matrix spike recovery criteria not met for this analyte
- U Not Detected
- NR Not Reported
- Q QC limit Exceeded



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 20012003\_2012005

**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:** Groundwater

**Collection date:** 12/01/2020 & 12/02/2020

**Analysis/Methods:**

Wet Chemistry:

Alkalinity M2320 B

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
2012003-01	A11-MW130A-201201	2012005-01	A11-MW007-201201-D
2012003-02	A11-MW006-201201	2012005-02	A11-MW007-201201
2012003-03	A11-MW005-201201	2012005-03	A11-MW004A-201201
2012003-04	A11-MW004B-201201	2012005-04	A11-MW003-201201
2012003-05	A11-MW001-201201	2012005-05	A11-MW002-201201
2012003-06	A11-FB001-201201		

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters (Alkalinity 2320B)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	N/A		
Laboratory Control Spike Duplicates RPD within limits?	N/A		
Laboratory Duplicate RPDs within limits?	Yes		
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b>	<b><u>Sample</u> A11-MW007-201201</b>	<b><u>Duplicate</u> A11-MW007-201201-D</b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
			Acceptable		

<b>MS/MSD</b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>LCS/LCSD</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20L009-DUP1	Acceptable			

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	N/A		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	Yes		
Was the ICAL criteria met?	N/A		
Was the CCV criteria met?	N/A		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

<b>Blanks</b>	<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20L009-BLK1	Nondetect			

Field Blank A11-FB001-201201	Concentration Nondetect	MDL / PQL	Qualifiers	Associated Samples										
Surrogates N/A	%R	Limit	Qualifiers	Associated Samples										
MS/MSD N/A	%R	Limits (%)	Qualifiers	Associated Samples										
LCS/LCSD B20L009-SRM1	%R Acceptable	Limits	Qualifiers	Associated Samples										
ICV N/A		%R Limits	Qualifiers	Associated Samples										
CCV N/A		%R Limits	Qualifiers	Associated Samples										
Tune N/A														
Internal Standards N/A	Area	Area Lower / Upper Limit	Qualifiers	Associated Samples										
<b>Representativeness:</b> <table border="1"> <tr> <td>Were sampling procedures and design criteria met?</td> <td>Yes No N/A</td> </tr> <tr> <td>Were holding times met?</td> <td>Yes</td> </tr> <tr> <td>Was preservation criteria met? (0° C - 6° C)</td> <td>Yes</td> </tr> <tr> <td>Were Chain-of-Custody records complete and provided in data package?</td> <td>No</td> </tr> <tr> <td colspan="2">Comments (note deviations): The cooler temperature was 0.3° C.</td> </tr> </table>					Were sampling procedures and design criteria met?	Yes No N/A	Were holding times met?	Yes	Was preservation criteria met? (0° C - 6° C)	Yes	Were Chain-of-Custody records complete and provided in data package?	No	Comments (note deviations): The cooler temperature was 0.3° C.	
Were sampling procedures and design criteria met?	Yes No N/A													
Were holding times met?	Yes													
Was preservation criteria met? (0° C - 6° C)	Yes													
Were Chain-of-Custody records complete and provided in data package?	No													
Comments (note deviations): The cooler temperature was 0.3° C.														
Preservation	Cooler Temperature (Degrees C) Acceptable	Preservation Criteria	Qualifier	Associated Samples										
Holding Times	Analyte	Days to Extraction Acceptable	HT Criteria	Qualifier Associated Samples										
<b>Comparability:</b> <table border="1"> <tr> <td>Were analytical procedures and methods followed as defined in the QAPP or field change documentation?</td> <td>Yes No N/A</td> </tr> <tr> <td>Comments (note deviations):</td> <td>Yes</td> </tr> </table>					Were analytical procedures and methods followed as defined in the QAPP or field change documentation?	Yes No N/A	Comments (note deviations):	Yes						
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?	Yes No N/A													
Comments (note deviations):	Yes													
<b>Completeness (90%):</b> <table border="1"> <tr> <td>Are all data in this SDG usable?</td> <td>Yes No N/A</td> </tr> <tr> <td>Comments (note deviations):</td> <td>Yes</td> </tr> </table>					Are all data in this SDG usable?	Yes No N/A	Comments (note deviations):	Yes						
Are all data in this SDG usable?	Yes No N/A													
Comments (note deviations):	Yes													
<b>Sensitivity:</b> <table border="1"> <tr> <td>Are MDLs present and reported?</td> <td>Yes No N/A</td> </tr> <tr> <td>Do the reporting limits meet project requirements?</td> <td>Yes</td> </tr> <tr> <td>Comments (note deviations):</td> <td>Yes</td> </tr> </table>					Are MDLs present and reported?	Yes No N/A	Do the reporting limits meet project requirements?	Yes	Comments (note deviations):	Yes				
Are MDLs present and reported?	Yes No N/A													
Do the reporting limits meet project requirements?	Yes													
Comments (note deviations):	Yes													
<b>Comment:</b> Data is usable as reported.														
Data Validator:	Kristine Molloy		Date: 1/20/2021/											
Data Reviewer:	Cherie Zakowski		Date: 1/23/2021											





**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination  
 Project Number: ILD981000417  
 Project Manager: Terese Van Donsel

**Reported:**  
 Dec-18-20 12:33

**Alkalinity by SM 2320B**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW130A-201201 (2012003-01) Matrix: Water Sampled: Dec-01-20 09:35 Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>350</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW006-201201 (2012003-02) Matrix: Water Sampled: Dec-01-20 12:11 Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>450</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW005-201201 (2012003-03) Matrix: Water Sampled: Dec-01-20 13:50 Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>380</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW004B-201201 (2012003-04) Matrix: Water Sampled: Dec-01-20 16:45 Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>340</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW001-201201 (2012003-05) Matrix: Water Sampled: Dec-01-20 05:31 Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>350</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-FB001-201201 (2012003-06) Matrix: Water Sampled: Dec-01-20 17:01 Received: Dec-02-20 11:34**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>U</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW007-201201-D (2012005-01) Matrix: Water Sampled: Dec-02-20 10:50 Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>540</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW007-201201 (2012005-02) Matrix: Water Sampled: Dec-02-20 10:50 Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>540</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20



## Environmental Protection Agency Region 5

### US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
 Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: SE Rockford GW Contamination

Project Number: ILD981000417

Project Manager: Terese Van Donsel

**Reported:**  
 Dec-18-20 12:33

### Alkalinity by SM 2320B

### US EPA Region 5 LSASD Analytical Services Branch

**A11-MW004A-201201 (2012005-03)**

**Matrix: Water**

**Sampled: Dec-02-20 15:00**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>330</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW003-201201 (2012005-04)**

**Matrix: Water**

**Sampled: Dec-02-20 08:45**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>390</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20

**A11-MW002-201201 (2012005-05)**

**Matrix: Water**

**Sampled: Dec-02-20 13:10**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Total Alkalinity</b>	<b>440</b>			20	mg CaCO3/L	1	B20L009	Dec-09-20	Dec-09-20





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Dec-18-20 12:33

**Notes and Definitions**

- \* This Quality Control measure meets the requirements of the CRL SOP for this analyte.
- U Not Detected
- NR Not Reported
- Q QC limit Exceeded

**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 2012003

**Laboratory:** ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:** Groundwater

**Collection date:** 12/01/20

**Analysis/Methods:** Wet Chemistry: Anions - EPA 300.0

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>
2012003-01	A11-MW130A-201201
2012003-02	A11-MW006-201201
2012003-03	A11-MW005-201201
2012003-04	A11-MW004B-201201
2012003-05	A11-MW001-201201
2012003-06	A11-FB001-201201

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Inorganic Superfund Methods Data Review (EPA January 2017).

**Wet Chemistry Parameters (Anions 300.0)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?			N/A
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)			N/A
Laboratory Control Spike Duplicates RPD within limits?			N/A
Laboratory Duplicate RPDs within limits?			Yes
<u>Comments (note deviations):</u>			

<b>Field Duplicates</b> N/A	<b><u>Sample</u></b>	<b><u>Duplicate</u></b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>MS/MSD</b> N/A	<b><u>%RPD</u></b>	<b><u>Limit</u></b>		<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>LCS/LCSD</b> N/A	<b><u>%RPD</u></b>	<b><u>Limits</u></b>		<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
<b>Laboratory Duplicate</b> E20L003-DUP1	<b><u>%RPD</u></b> Acceptable	<b><u>Limits</u></b>		<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	Yes		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

<b>Blanks</b>	<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
E20L003-BLK1				
Nitrogen, Nitrate	Nondetect			
Sulfate	Nondetect			



<b>ICB/CCB</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
ICB	Nitrogen, Nitrate Sulfate	Nondetect 0.03	0.10 / 0.12					
CCB1	Nitrogen, Nitrate Sulfate	Nondetect 0.03	0.10 / 0.12	None	Sample results nondetect or > RL			
<b>Field Blank</b>		<b><u>Concentration</u></b>	<b><u>MDL / PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
A11-FB001-201201		Nondetect						
<b>Surrogates</b>		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
N/A								
<b>MS/MSD</b>		<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
E20L003-MS1								
Nitrogen, Nitrate Sulfate		Acceptable 68%	80-120 80-120	J- / UJ	All samples			
<b>LCS/LCSD</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
E20L003-BS1								
Nitrogen, Nitrate Sulfate		Acceptable Acceptable	90-110 90-110					
<b>ICV</b>			<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
	Nitrogen, Nitrate Sulfate		Acceptable Acceptable					
<b>CCV</b>			<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
	Nitrogen, Nitrate Sulfate		Acceptable Acceptable					
<b>MRL Check</b>			<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>		
B20L003-MRL1								
Nitrogen, Nitrate Sulfate			Acceptable Acceptable					
<b>Tune</b>								
N/A								
<b>Internal Standards</b>		<b><u>Area</u></b>	<b><u>Area Lower / Upper Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>			
N/A								
<b>Representativeness:</b>						<b><u>Yes</u></b>	<b><u>No</u></b>	<b><u>N/A</u></b>
Were sampling procedures and design criteria met?							Yes	
Were holding times met?							Yes	
Was preservation criteria met? (0° C - 6° C)							Yes	
Were Chain-of-Custody records complete and provided in data package?							No	
<u>Comments (note deviations):</u> The cooler temperature was 0.3 ° C.								
<b>Preservation</b>		<b><u>Cooler Temperature (Degrees C)</u></b>	<b><u>Preservation Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>			
		Acceptable						
<b>Holding Times</b>		<b><u>Analyte</u></b>	<b><u>Days to Extraction</u></b>	<b><u>HT Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>		
			Acceptable					

**Comparability:**

Were analytical procedures and methods followed as defined in the QAPP or field change documentation?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Completeness (90%):**

Are all data in this SDG usable?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Sensitivity:**

Are MDLs present and reported?

Do the reporting limits meet project requirements?

Comments (note deviations):

**Yes No N/A**

**Yes**

**Yes**

**Comment:**

Data is usable with appropriate qualifiers applied.

Data Validator:

*Kristine Molloy*

Date: 1/19/2021

Data Reviewer:

Cherie Zakowski

Date: 1/25/2021





## Environmental Protection Agency Region 5

### US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone: (312) 353-8370 Fax: (312) 886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Dec-07-20 16:52

### Anions by Ion Chromatography, EPA 300.0 (modified) US EPA Region 5 LSASD Analytical Services Branch

**A11-MW007-201201-D (2012005-01)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	2.56			0.12	mg/L	1	B20L006	Dec-03-20	Dec-03-20
Nitrate - NO <sub>3</sub>	U			0.12	"	"	"	"	"

**A11-MW007-201201 (2012005-02)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	2.45			0.12	mg/L	1	B20L006	Dec-03-20	Dec-03-20
Nitrate - NO <sub>3</sub>	U			0.12	"	"	"	"	"

**A11-MW004A-201201 (2012005-03)**

**Matrix: Water**

**Sampled: Dec-02-20 15:00**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	42.9			0.12	mg/L	1	B20L006	Dec-03-20	Dec-03-20
Nitrate - NO <sub>3</sub>	1.66			0.12	"	"	"	"	"

**A11-MW003-201201 (2012005-04)**

**Matrix: Water**

**Sampled: Dec-02-20 08:45**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	8.52			0.12	mg/L	1	B20L006	Dec-03-20	Dec-03-20
Nitrate - NO <sub>3</sub>	U			0.12	"	"	"	"	"

**A11-MW002-201201 (2012005-05)**

**Matrix: Water**

**Sampled: Dec-02-20 13:10**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Sulfate as SO <sub>4</sub>	1.09			0.12	mg/L	1	B20L006	Dec-03-20	Dec-03-20
Nitrate - NO <sub>3</sub>	U			0.12	"	"	"	"	"



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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Dec-07-20 16:52

**Notes and Definitions**

U Not Detected  
NR Not Reported  
Q QC limit Exceeded



**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:**

2012003, 2012005

**Laboratory:**

ESAT - US EPA Region 5 LSASD Analytical Services Branch

**Matrix:**

Groundwater

**Collection date:**

12/01/2020 & 12/02/2020

**Analysis/Methods:**

Volatile Organic Compounds (VOCs) 8260

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
2012003-01RE1	A11-MW130A-201201	2012005-01RE1	A11-MW007-201201-D
2012003-02RE1	A11-MW006-201201	2012005-02RE1	A11-MW007-201201
2012003-03RE1	A11-MW005-201201	2012005-03RE2	A11-MW004A-201201
2012003-04RE1	A11-MW004B-201201	2012005-04RE1	A11-MW003-201201
2012003-05RE1	A11-MW001-201201	2012005-05RE2	A11-MW002-201201
2012003-06	A11-FB001-201201	2012005-06RE1	A11-TB002-201201
2012003-07	A11-TB001-201201		

Data validation was performed in accordance with the specific analytical method and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Volatile Organic Compounds 8260 / 1,4-Dioxane 8000D**

**Precision:**

Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?

**Yes No N/A**

**No**

Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)

**Yes**

Laboratory Control Spike Duplicates RPD within limits?

**No**

Laboratory Duplicate RPDs within limits?

**N/A**

Comments (note deviations):

<b>Field Duplicates</b>	<b><u>Sample</u></b>	<b><u>Duplicate</u></b>	<b><u>%RPD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
	<b>A11-MW007-201201</b>	<b>A11-MW007-201201-D</b>			
Isopropylbenzene	109	486	127%	J	<b>A11-MW007-201201 &amp; A11-MW007-201201-D</b>
n-Propylbenzene	104	454	125%	J	
sec-Butylbenzene	17.5	68.4	74%	J	
1,3,5-Trimethylbenzene	14.4	56.7	NC	J*	
Benzene	10 U	44.3	NC	J / UJ*	
Naphthalene	34	97.3	NC	J*	Sample results < 5xs RL; ABS Diff. < RL
n-Butylbenzene	19.9	66.9	NC	J*	
1,2,4-Trimethylbenzene	131	169	NC	None	

\* Sample results < 5xs RL; ABS Diff. > RL

<b><u>MS/MSD</u></b>	<b><u>%RPD</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20L005-MS1 / MSD1 (2012003-03RE1)	Acceptable			

<b><u>LCS/LCSD</u></b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
B20L004-BS1 / BSD1 Acetone	23.5	20%	J**	2012003-06, 2012003-07
B20L005-BS1 / BSD1 2,2-Dichloropropane	23.7	20%	J**	2012003-01RE1 through 2012003-05RE1
B20L008-BS1 / BSD1	Acceptable			
B20L008-BS2 / BSD2 2,2-Dichloropropane	57.8	20%	J**	2012005-03RE2, 2012005-05RE2

\*\*Qualification required for detected results only - associated results nondetect - no qualification required

<b>Laboratory Duplicate</b>	<b><u>%RPD</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
N/A				

Accuracy:	Yes	No	N/A
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	No		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	Yes		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	Yes		
Were the Surrogate % recoveries within laboratory determined control limits?	Yes		
Were the Internal Standard areas within ± 50 - 150%?	Yes		
Comments (note deviations):			

Blanks	Concentration	MDL /PQL	Qualifiers	Associated Samples
B20L004-BLK1	Nondetect			
B20L004-BLK2	Nondetect			
B20L005-BLK1	Nondetect			
B20L005-BLK2	Nondetect			
B20L008-BLK1	Nondetect			
B20L008-BLK2	Nondetect			

Field Blank	Concentration	MDL / PQL	Qualifiers	Associated Samples
A11-FB001-201201	Nondetect			
A11-TB001-201201	Nondetect			
A11-TB001-201201	Nondetect			

Surrogates	%R	Limit	Qualifiers	Associated Samples
	Acceptable			

MS/MSD	%R	Limits (%)	Qualifiers	Associated Samples
B20L005-MS1 / MSD1 (2012003-03RE1)	Acceptable			

LCS/LCSD		%R	Limits	Qualifiers	Associated Samples
B20L004-BS1 / BSD1		Acceptable			
B20L005-BS1 / BSD1		Acceptable			
B20L005-BS2		Acceptable			
B20L008-BS1 / BSD1		Acceptable			
B20L008-BS2/ BSD2	2,2-Dichloropropane	71.5 / 39.5	70-130	J / UJ	2012005-03RE2, 2012005-05RE2

ICAL	RRF	%RSD	Limits	Qualifiers	Associated Samples
12/1/2020 11:44	Acceptable	Acceptable			

ICV / CCV	RRF	%D	Limits	Qualifiers	Associated Samples
ICV					
12/1/2020 3:28	Acceptable	Acceptable			
CCV					
12/2/2020 13:34	Acceptable	Acceptable			
12/2/2020 18:26	Acceptable	Acceptable			
12/3/2020 9:03	Acceptable	Acceptable			
12/3/2020 13:44	Acceptable	Acceptable			
12/3/2020 19:35	Acceptable	Acceptable			
12/4/2020 12:40	Acceptable	Acceptable			
12/4/2020 18:54	Acceptable	Acceptable			
12/5/2020 1:58	Acceptable	Acceptable			
12/5/2020 9:28	Acceptable	Acceptable			



<b>MRL Check</b>	<u>%R</u>	<u>Limits</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
B20L005-MRL1	Acceptable			

<b>Tune</b>
Acceptable

<b>Internal Standards</b>	<u>Area</u>	<u>Area Lower / Upper Limit</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
		Acceptable		

<b>Representativeness:</b>	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Were sampling procedures and design criteria met?	Yes		
Were holding times met?	Yes		
Was preservation criteria met? (0° C - 6° C)	Yes		
Were Chain-of-Custody records complete and provided in data package?	Yes		
<u>Comments (note deviations):</u> The cooler temperature was 0.30 ° C.			

<b>Preservation</b>	<u>Cooler Temperature (Degrees C)</u>	<u>Preservation Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
	Acceptable			

<b>Holding Times</b>	<u>Analyte</u>	<u>Days to Extraction</u>	<u>HT Criteria</u>	<u>Qualifier</u>	<u>Associated Samples</u>
		Acceptable			

<b>Comparability:</b>	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?	Yes		
<u>Comments (note deviations):</u>			

<b>Completeness (90%):</b>	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Are all data in this SDG usable?	Yes		
<u>Comments (note deviations):</u>			

<b>Sensitivity:</b>	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Are MDLs present and reported?	Yes		
Do the reporting limits meet project requirements?	Yes		
<u>Comments (note deviations):</u>			

**Comment:**

As stated in the case narrative, sample 2012003-03 vial B was not acid preserved, but was analyzed within 7 days of sampling, All other samples were properly preserved (acidification in field) and met hold time (14 days) criteria.

As stated in the case narrative, samples were first screened at a 50x dilution and some re-injections were required. Each analyte is reported at the lowest dilution factor for which the analyte concentration remained within calibration range.

As stated in the case narrative, no matrix spike was analyzed for the samples associated with batch B2L008 due to insufficient number of vials.

Case narrative indicates co-elution affected the concentration of n-butylbenzene, n-Butylbenzene has been flagged as an estimated concentration (J) when detected above the reporting limit.

Data is usable with appropriate qualifiers applied.

Data Validator:	<u>Kristine Molloy</u>	Date:	<u>5/3/2021</u>
Data Reviewer:	<u>Cherie Zakowski</u>	Date:	<u>5/6/2021</u>



Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW130A-201201 (2012003-01RE1)

Matrix: Water

Sampled: Dec-01-20 09:35

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	3.77			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	3.51			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"





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A11-MW130A-201201 (2012003-01RE1)

Matrix: Water

Sampled: Dec-01-20 09:35

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Chlorobenzene	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.57			95.1%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			101%		84-122	"	"	"
Toluene-d8	9.66			96.6%		88-108	"	"	"
4-Bromofluorobenzene	9.86			98.6%		84-108	"	"	"



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Project Number: ILD981000417  
Project Manager: Terese Van Donsel

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Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW006-201201 (2012003-02RE1)

Matrix: Water

Sampled: Dec-01-20 12:11

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	2.82			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Matrix: Water

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Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.2			101%		73-124	"	"	"
1,2-Dichloroethane-d4	10.6			105%		84-122	"	"	"
Toluene-d8	9.95			99.5%		88-108	"	"	"
4-Bromofluorobenzene	10.2			102%		84-108	"	"	"



**Environmental Protection Agency Region 5**  
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**Reported:**  
Jan-15-21 13:14

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-MW005-201201 (2012003-03RE1)

Matrix: Water

Sampled: Dec-01-20 13:50

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	7.01			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	4.90			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.75			96.9%		73-124	"	"	"
1,2-Dichloroethane-d4	10.1			100%		84-122	"	"	"
Toluene-d8	10.1			101%		88-108	"	"	"
4-Bromofluorobenzene	9.55			95.5%		84-108	"	"	"



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A11-MW004B-201201 (2012003-04RE1)

Matrix: Water

Sampled: Dec-01-20 16:45

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	5.67			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	5.61			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW004B-201201 (2012003-04RE1)

Matrix: Water

Sampled: Dec-01-20 16:45

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.88			98.2%		73-124	"	"	"
1,2-Dichloroethane-d4	10.2			101%		84-122	"	"	"
Toluene-d8	9.87			98.7%		88-108	"	"	"
4-Bromofluorobenzene	9.68			96.8%		84-108	"	"	"



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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW001-201201 (2012003-05RE1)

Matrix: Water

Sampled: Dec-01-20 05:31

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	4.94			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	9.02			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	2.15			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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77 West Jackson Boulevard  
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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW001-201201 (2012003-05RE1)

Matrix: Water

Sampled: Dec-01-20 05:31

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20L005	Dec-02-20	Dec-03-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.71			96.5%		73-124	"	"	"
1,2-Dichloroethane-d4	10.1			100%		84-122	"	"	"
Toluene-d8	10.0			100%		88-108	"	"	"
4-Bromofluorobenzene	9.69			96.9%		84-108	"	"	"



## Environmental Protection Agency Region 5

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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

### Volatiles by GC/MS, EPA 8260C (modified) US EPA Region 5 LSASD Analytical Services Branch

A11-FB001-201201 (2012003-06)

Matrix: Water

Sampled: Dec-01-20 17:01

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L004	Dec-02-20	Dec-02-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-FB001-201201 (2012003-06)

Matrix: Water

Sampled: Dec-01-20 17:01

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20L004	Dec-02-20	Dec-02-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.83			97.8%		73-124	"	"	"
1,2-Dichloroethane-d4	9.94			98.7%		84-122	"	"	"
Toluene-d8	9.80			98.0%		88-108	"	"	"
4-Bromofluorobenzene	9.99			99.9%		84-108	"	"	"



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Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB001-201201 (2012003-07)

Matrix: Water

Sampled: Dec-01-20 09:00

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L004	Dec-02-20	Dec-02-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"
Chloroform	U			2.00	"	"	"	"	"
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"





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Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB001-201201 (2012003-07)

Matrix: Water

Sampled: Dec-01-20 09:00

Received: Dec-02-20 11:34

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,1,1,2-Tetrachloroethane	U			2.00	ug/L	1	B20L004	Dec-02-20	Dec-02-20
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"
sec-Butylbenzene	U			2.00	"	"	"	"	"
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.1			100%		73-124	"	"	"
1,2-Dichloroethane-d4	10.4			103%		84-122	"	"	"
Toluene-d8	9.86			98.6%		88-108	"	"	"
4-Bromofluorobenzene	9.87			98.7%		84-108	"	"	"



**Environmental Protection Agency Region 5**  
**US EPA Region 5 LSASD Analytical Services Branch**

536 South Clark Street, Chicago, IL 60605  
Phone: (312) 353-8370 Fax: (312) 886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

**Reported:**  
Jan-15-21 13:14

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW007-201201-D (2012005-01)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Ethylbenzene</b>	<b>3660</b>			100	ug/L	50	B20L005	Dec-02-20	Dec-03-20
<b>m+p-Xylene</b>	<b>8100</b>			200	"	"	"	"	"
<b>1,2,4-Trimethylbenzene</b>	<b>169</b>			100	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
<i>Dibromofluoromethane</i>	9.76			97.0%		73-124	"	"	"
<i>1,2-Dichloroethane-d4</i>	10.1			99.8%		84-122	"	"	"
<i>Toluene-d8</i>	9.78			97.8%		88-108	"	"	"
<i>4-Bromofluorobenzene</i>	9.83			98.3%		84-108	"	"	"

**A11-MW007-201201-D (2012005-01RE1)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Dichlorodifluoromethane</b>	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
<b>Chloromethane</b>	U			10.0	"	"	"	"	"
<b>Vinyl chloride</b>	U			10.0	"	"	"	"	"
<b>Bromomethane</b>	U			10.0	"	"	"	"	"
<b>Chloroethane</b>	U			10.0	"	"	"	"	"
<b>Trichlorofluoromethane</b>	U			10.0	"	"	"	"	"
<b>1,1-Dichloroethene</b>	U			10.0	"	"	"	"	"
<b>Acetone</b>	U			62.5	"	"	"	"	"
<b>Carbon disulfide</b>	U			10.0	"	"	"	"	"
<b>Methylene chloride</b>	U			10.0	"	"	"	"	"
<b>trans-1,2-Dichloroethene</b>	U			10.0	"	"	"	"	"
<b>1,1-Dichloroethane</b>	U			10.0	"	"	"	"	"
<b>2,2-Dichloropropane</b>	U			10.0	"	"	"	"	"
<b>cis-1,2-Dichloroethene</b>	U			10.0	"	"	"	"	"
<b>2-Butanone</b>	U			62.5	"	"	"	"	"
<b>Bromochloromethane</b>	U			10.0	"	"	"	"	"
<b>Chloroform</b>	U			10.0	"	"	"	"	"
<b>1,1,1-Trichloroethane</b>	U			10.0	"	"	"	"	"
<b>Carbon tetrachloride</b>	U			10.0	"	"	"	"	"
<b>1,1-Dichloropropene</b>	U			10.0	"	"	"	"	"
<b>Benzene</b>	<b>44.3</b>			10.0	"	"	"	"	"
<b>1,2-Dichloroethane</b>	U			10.0	"	"	"	"	"
<b>Trichloroethene</b>	U			10.0	"	"	"	"	"
<b>1,2-Dichloropropane</b>	U			10.0	"	"	"	"	"





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 Jan-15-21 13:14

**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-MW007-201201-D (2012005-01RE1)

Matrix: Water

Sampled: Dec-02-20 10:50

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dibromomethane	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
Bromodichloromethane	U			10.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			10.0	"	"	"	"	"
4-Methyl-2-pentanone	U			25.0	"	"	"	"	"
Toluene	U			10.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			10.0	"	"	"	"	"
1,1,2-Trichloroethane	U			10.0	"	"	"	"	"
Tetrachloroethene	U			10.0	"	"	"	"	"
1,3-Dichloropropane	U			10.0	"	"	"	"	"
2-Hexanone	U			25.0	"	"	"	"	"
Dibromochloromethane	U			10.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			10.0	"	"	"	"	"
Chlorobenzene	U			10.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			10.0	"	"	"	"	"
o-Xylene	U			10.0	"	"	"	"	"
Styrene	U			10.0	"	"	"	"	"
Bromoform	U			10.0	"	"	"	"	"
Isopropylbenzene	486			10.0	"	"	"	"	"
Bromobenzene	U			10.0	"	"	"	"	"
1,2,3-Trichloropropane	U			10.0	"	"	"	"	"
n-Propylbenzene	454			10.0	"	"	"	"	"
2-Chlorotoluene	U			10.0	"	"	"	"	"
1,3,5-Trimethylbenzene	56.7			10.0	"	"	"	"	"
4-Chlorotoluene	U			10.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			10.0	"	"	"	"	"
tert-Butylbenzene	U			10.0	"	"	"	"	"
sec-Butylbenzene	68.4			10.0	"	"	"	"	"
1,3-Dichlorobenzene	U			10.0	"	"	"	"	"
p-Isopropyltoluene	U			10.0	"	"	"	"	"
1,4-Dichlorobenzene	U			10.0	"	"	"	"	"
1,2-Dichlorobenzene	U			10.0	"	"	"	"	"
n-Butylbenzene	66.9	CustomFlag, J		10.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			10.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			10.0	"	"	"	"	"
Hexachlorobutadiene	U			10.0	"	"	"	"	"
Naphthalene	97.3			10.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			10.0	"	"	"	"	"



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**US EPA Region 5 LSASD Analytical Services Branch**

**A11-MW007-201201-D (2012005-01RE1)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.97			99.1%		73-124	B20L008	Dec-04-20	Dec-04-20
1,2-Dichloroethane-d4	10.4			104%		84-122	"	"	"
Toluene-d8	10.3			103%		88-108	"	"	"
4-Bromofluorobenzene	9.19			91.9%		84-108	"	"	"

**A11-MW007-201201 (2012005-02)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethylbenzene	3300			100	ug/L	50	B20L005	Dec-02-20	Dec-03-20
m+p-Xylene	7390			200	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.2			102%		73-124	"	"	"
1,2-Dichloroethane-d4	10.3			102%		84-122	"	"	"
Toluene-d8	10.1			101%		88-108	"	"	"
4-Bromofluorobenzene	9.78			97.8%		84-108	"	"	"

**A11-MW007-201201 (2012005-02RE1)**

**Matrix: Water**

**Sampled: Dec-02-20 10:50**

**Received: Dec-03-20 10:57**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
Chloromethane	U			10.0	"	"	"	"	"
Vinyl chloride	U			10.0	"	"	"	"	"
Bromomethane	U			10.0	"	"	"	"	"
Chloroethane	U			10.0	"	"	"	"	"
Trichlorofluoromethane	U			10.0	"	"	"	"	"
1,1-Dichloroethene	U			10.0	"	"	"	"	"
Acetone	U			62.5	"	"	"	"	"
Carbon disulfide	U			10.0	"	"	"	"	"
Methylene chloride	U			10.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			10.0	"	"	"	"	"
1,1-Dichloroethane	U			10.0	"	"	"	"	"
2,2-Dichloropropane	U			10.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			10.0	"	"	"	"	"
2-Butanone	U			62.5	"	"	"	"	"





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A11-MW007-201201 (2012005-02RE1)

Matrix: Water

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Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Bromochloromethane	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
Chloroform	U			10.0	"	"	"	"	"
1,1,1-Trichloroethane	U			10.0	"	"	"	"	"
Carbon tetrachloride	U			10.0	"	"	"	"	"
1,1-Dichloropropene	U			10.0	"	"	"	"	"
Benzene	U			10.0	"	"	"	"	"
1,2-Dichloroethane	U			10.0	"	"	"	"	"
Trichloroethene	U			10.0	"	"	"	"	"
1,2-Dichloropropane	U			10.0	"	"	"	"	"
Dibromomethane	U			10.0	"	"	"	"	"
Bromodichloromethane	U			10.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			10.0	"	"	"	"	"
4-Methyl-2-pentanone	U			25.0	"	"	"	"	"
Toluene	U			10.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			10.0	"	"	"	"	"
1,1,2-Trichloroethane	U			10.0	"	"	"	"	"
Tetrachloroethene	U			10.0	"	"	"	"	"
1,3-Dichloropropane	U			10.0	"	"	"	"	"
2-Hexanone	U			25.0	"	"	"	"	"
Dibromochloromethane	U			10.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			10.0	"	"	"	"	"
Chlorobenzene	U			10.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			10.0	"	"	"	"	"
o-Xylene	U			10.0	"	"	"	"	"
Styrene	U			10.0	"	"	"	"	"
Bromoform	U			10.0	"	"	"	"	"
Isopropylbenzene	109			10.0	"	"	"	"	"
Bromobenzene	U			10.0	"	"	"	"	"
1,2,3-Trichloropropane	U			10.0	"	"	"	"	"
n-Propylbenzene	104			10.0	"	"	"	"	"
2-Chlorotoluene	U			10.0	"	"	"	"	"
1,3,5-Trimethylbenzene	14.4			10.0	"	"	"	"	"
4-Chlorotoluene	U			10.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			10.0	"	"	"	"	"
tert-Butylbenzene	U			10.0	"	"	"	"	"
1,2,4-Trimethylbenzene	131			10.0	"	"	"	"	"
sec-Butylbenzene	17.5			10.0	"	"	"	"	"



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A11-MW007-201201 (2012005-02RE1)

Matrix: Water

Sampled: Dec-02-20 10:50

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,3-Dichlorobenzene	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
p-Isopropyltoluene	U			10.0	"	"	"	"	"
1,4-Dichlorobenzene	U			10.0	"	"	"	"	"
1,2-Dichlorobenzene	U			10.0	"	"	"	"	"
n-Butylbenzene	19.9	B, CustomFlag, J		10.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			10.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			10.0	"	"	"	"	"
Hexachlorobutadiene	U			10.0	"	"	"	"	"
Naphthalene	34.0	B		10.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			10.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.0			99.6%		73-124	"	"	"
1,2-Dichloroethane-d4	10.3			102%		84-122	"	"	"
Toluene-d8	9.82			98.2%		88-108	"	"	"
4-Bromofluorobenzene	9.25			92.5%		84-108	"	"	"

A11-MW004A-201201 (2012005-03RE1)

Matrix: Water

Sampled: Dec-02-20 15:00

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Toluene	34200			1250	ug/L	625	B20L008	Dec-04-20	Dec-04-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.83			97.7%		73-124	"	"	"
1,2-Dichloroethane-d4	10.6			105%		84-122	"	"	"
Toluene-d8	10.1			101%		88-108	"	"	"
4-Bromofluorobenzene	9.78			97.8%		84-108	"	"	"

A11-MW004A-201201 (2012005-03RE2)

Matrix: Water

Sampled: Dec-02-20 15:00

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			50.0	ug/L	25	B20L008	Dec-04-20	Dec-05-20
Chloromethane	U			50.0	"	"	"	"	"
Vinyl chloride	U			50.0	"	"	"	"	"
Bromomethane	U			50.0	"	"	"	"	"
Chloroethane	U			50.0	"	"	"	"	"





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**Volatiles by GC/MS, EPA 8260C (modified)**  
**US EPA Region 5 LSASD Analytical Services Branch**

A11-MW004A-201201 (2012005-03RE2)

Matrix: Water

Sampled: Dec-02-20 15:00

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Trichlorofluoromethane	U			50.0	ug/L	25	B20L008	Dec-04-20	Dec-05-20
1,1-Dichloroethene	U			50.0	"	"	"	"	"
Acetone	U			312	"	"	"	"	"
Carbon disulfide	U			50.0	"	"	"	"	"
Methylene chloride	U			50.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			50.0	"	"	"	"	"
1,1-Dichloroethane	U			50.0	"	"	"	"	"
2,2-Dichloropropane	U	(LCS), J		50.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			50.0	"	"	"	"	"
2-Butanone	U			312	"	"	"	"	"
Bromochloromethane	U			50.0	"	"	"	"	"
Chloroform	U			50.0	"	"	"	"	"
1,1,1-Trichloroethane	U			50.0	"	"	"	"	"
Carbon tetrachloride	U			50.0	"	"	"	"	"
1,1-Dichloropropene	U			50.0	"	"	"	"	"
Benzene	U			50.0	"	"	"	"	"
1,2-Dichloroethane	U			50.0	"	"	"	"	"
Trichloroethene	U			50.0	"	"	"	"	"
1,2-Dichloropropane	U			50.0	"	"	"	"	"
Dibromomethane	U			50.0	"	"	"	"	"
Bromodichloromethane	U			50.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			50.0	"	"	"	"	"
4-Methyl-2-pentanone	U			125	"	"	"	"	"
trans-1,3-Dichloropropene	U			50.0	"	"	"	"	"
1,1,2-Trichloroethane	U			50.0	"	"	"	"	"
Tetrachloroethene	U			50.0	"	"	"	"	"
1,3-Dichloropropane	U			50.0	"	"	"	"	"
2-Hexanone	U			125	"	"	"	"	"
Dibromochloromethane	U			50.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			50.0	"	"	"	"	"
Chlorobenzene	U			50.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			50.0	"	"	"	"	"
Ethylbenzene	331			50.0	"	"	"	"	"
m+p-Xylene	489			100	"	"	"	"	"
o-Xylene	52.2			50.0	"	"	"	"	"
Styrene	U			50.0	"	"	"	"	"
Bromoform	U			50.0	"	"	"	"	"



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A11-MW004A-201201 (2012005-03RE2)

Matrix: Water

Sampled: Dec-02-20 15:00

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Isopropylbenzene	U			50.0	ug/L	25	B20L008	Dec-04-20	Dec-05-20
Bromobenzene	U			50.0	"	"	"	"	"
1,2,3-Trichloropropane	U			50.0	"	"	"	"	"
n-Propylbenzene	U			50.0	"	"	"	"	"
2-Chlorotoluene	U			50.0	"	"	"	"	"
1,3,5-Trimethylbenzene	U			50.0	"	"	"	"	"
4-Chlorotoluene	U			50.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			50.0	"	"	"	"	"
tert-Butylbenzene	U			50.0	"	"	"	"	"
1,2,4-Trimethylbenzene	U			50.0	"	"	"	"	"
sec-Butylbenzene	U			50.0	"	"	"	"	"
1,3-Dichlorobenzene	U			50.0	"	"	"	"	"
p-Isopropyltoluene	U			50.0	"	"	"	"	"
1,4-Dichlorobenzene	U			50.0	"	"	"	"	"
1,2-Dichlorobenzene	U			50.0	"	"	"	"	"
n-Butylbenzene	U			50.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			50.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			50.0	"	"	"	"	"
Hexachlorobutadiene	U			50.0	"	"	"	"	"
Naphthalene	U			50.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			50.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.97			99.1%		73-124	"	"	"
1,2-Dichloroethane-d4	10.5			104%		84-122	"	"	"
Toluene-d8	10.3			103%		88-108	"	"	"
4-Bromofluorobenzene	9.76			97.6%		84-108	"	"	"





Environmental Protection Agency Region 5  
US EPA Region 5 LSASD Analytical Services Branch

536 South Clark Street, Chicago, IL 60605  
Phone:(312)353-8370 Fax:(312)886-2591

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: SE Rockford GW Contamination  
Project Number: ILD981000417  
Project Manager: Terese Van Donsel

Reported:  
Jan-15-21 13:14

Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW003-201201 (2012005-04)

Matrix: Water

Sampled: Dec-02-20 08:45

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
m+p-Xylene	6310			200	ug/L	50	B20L005	Dec-02-20	Dec-03-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.86			98.0%		73-124	"	"	"
1,2-Dichloroethane-d4	10.4			103%		84-122	"	"	"
Toluene-d8	9.68			96.8%		88-108	"	"	"
4-Bromofluorobenzene	9.52			95.2%		84-108	"	"	"

A11-MW003-201201 (2012005-04RE1)

Matrix: Water

Sampled: Dec-02-20 08:45

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
Chloromethane	U			10.0	"	"	"	"	"
Vinyl chloride	U			10.0	"	"	"	"	"
Bromomethane	U			10.0	"	"	"	"	"
Chloroethane	U			10.0	"	"	"	"	"
Trichlorofluoromethane	U			10.0	"	"	"	"	"
1,1-Dichloroethene	U			10.0	"	"	"	"	"
Acetone	U			62.5	"	"	"	"	"
Carbon disulfide	U			10.0	"	"	"	"	"
Methylene chloride	U			10.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			10.0	"	"	"	"	"
1,1-Dichloroethane	U			10.0	"	"	"	"	"
2,2-Dichloropropane	U			10.0	"	"	"	"	"
cis-1,2-Dichloroethene	U			10.0	"	"	"	"	"
2-Butanone	U			62.5	"	"	"	"	"
Bromochloromethane	U			10.0	"	"	"	"	"
Chloroform	U			10.0	"	"	"	"	"
1,1,1-Trichloroethane	U			10.0	"	"	"	"	"
Carbon tetrachloride	U			10.0	"	"	"	"	"
1,1-Dichloropropene	U			10.0	"	"	"	"	"
Benzene	U			10.0	"	"	"	"	"
1,2-Dichloroethane	U			10.0	"	"	"	"	"
Trichloroethene	U			10.0	"	"	"	"	"
1,2-Dichloropropane	U			10.0	"	"	"	"	"
Dibromomethane	U			10.0	"	"	"	"	"
Bromodichloromethane	U			10.0	"	"	"	"	"



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Reported:  
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Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW003-201201 (2012005-04RE1)

Matrix: Water

Sampled: Dec-02-20 08:45

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
cis-1,3-Dichloropropene	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
4-Methyl-2-pentanone	U			25.0	"	"	"	"	"
Toluene	U			10.0	"	"	"	"	"
trans-1,3-Dichloropropene	U			10.0	"	"	"	"	"
1,1,2-Trichloroethane	U			10.0	"	"	"	"	"
Tetrachloroethene	U			10.0	"	"	"	"	"
1,3-Dichloropropane	U			10.0	"	"	"	"	"
2-Hexanone	U			25.0	"	"	"	"	"
Dibromochloromethane	U			10.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			10.0	"	"	"	"	"
Chlorobenzene	U			10.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			10.0	"	"	"	"	"
Ethylbenzene	256			10.0	"	"	"	"	"
o-Xylene	U			10.0	"	"	"	"	"
Styrene	U			10.0	"	"	"	"	"
Bromoform	U			10.0	"	"	"	"	"
Isopropylbenzene	38.5			10.0	"	"	"	"	"
Bromobenzene	U			10.0	"	"	"	"	"
1,2,3-Trichloropropane	U			10.0	"	"	"	"	"
n-Propylbenzene	37.3			10.0	"	"	"	"	"
2-Chlorotoluene	U			10.0	"	"	"	"	"
1,3,5-Trimethylbenzene	55.0			10.0	"	"	"	"	"
4-Chlorotoluene	U			10.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			10.0	"	"	"	"	"
tert-Butylbenzene	U			10.0	"	"	"	"	"
1,2,4-Trimethylbenzene	178			10.0	"	"	"	"	"
sec-Butylbenzene	15.0			10.0	"	"	"	"	"
1,3-Dichlorobenzene	U			10.0	"	"	"	"	"
p-Isopropyltoluene	U			10.0	"	"	"	"	"
1,4-Dichlorobenzene	U			10.0	"	"	"	"	"
1,2-Dichlorobenzene	U			10.0	"	"	"	"	"
n-Butylbenzene	13.1	B, CustomFlag, J		10.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			10.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			10.0	"	"	"	"	"
Hexachlorobutadiene	U			10.0	"	"	"	"	"
Naphthalene	16.6	B		10.0	"	"	"	"	"





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Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW003-201201 (2012005-04RE1)

Matrix: Water

Sampled: Dec-02-20 08:45

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2,3-Trichlorobenzene	U			10.0	ug/L	5	B20L008	Dec-04-20	Dec-04-20
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.81			97.5%		73-124	"	"	"
1,2-Dichloroethane-d4	10.9			108%		84-122	"	"	"
Toluene-d8	9.91			99.1%		88-108	"	"	"
4-Bromofluorobenzene	9.79			97.9%		84-108	"	"	"

A11-MW002-201201 (2012005-05RE1)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Toluene	33200			1250	ug/L	625	B20L008	Dec-04-20	Dec-04-20
Ethylbenzene	10200			1250	"	"	"	"	"
m+p-Xylene	31900			2500	"	"	"	"	"
o-Xylene	6140			1250	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.1			100%		73-124	"	"	"
1,2-Dichloroethane-d4	10.7			106%		84-122	"	"	"
Toluene-d8	9.89			98.9%		88-108	"	"	"
4-Bromofluorobenzene	9.65			96.5%		84-108	"	"	"

A11-MW002-201201 (2012005-05RE2)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			50.0	ug/L	25	B20L008	Dec-04-20	Dec-05-20
Chloromethane	U			50.0	"	"	"	"	"
Vinyl chloride	U			50.0	"	"	"	"	"
Bromomethane	U			50.0	"	"	"	"	"
Chloroethane	U			50.0	"	"	"	"	"
Trichlorofluoromethane	U			50.0	"	"	"	"	"
1,1-Dichloroethene	U			50.0	"	"	"	"	"
Acetone	U			312	"	"	"	"	"
Carbon disulfide	U			50.0	"	"	"	"	"
Methylene chloride	U			50.0	"	"	"	"	"
trans-1,2-Dichloroethene	U			50.0	"	"	"	"	"
1,1-Dichloroethane	U			50.0	"	"	"	"	"



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77 West Jackson Boulevard  
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**Reported:**  
Jan-15-21 13:14

### Volatiles by GC/MS, EPA 8260C (modified) US EPA Region 5 LSASD Analytical Services Branch

A11-MW002-201201 (2012005-05RE2)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
2,2-Dichloropropane	U	(LCS), J		50.0	ug/L	25	B20L008	Dec-04-20	Dec-05-20
cis-1,2-Dichloroethene	U			50.0	"	"	"	"	"
2-Butanone	U			312	"	"	"	"	"
Bromochloromethane	U			50.0	"	"	"	"	"
Chloroform	U			50.0	"	"	"	"	"
1,1,1-Trichloroethane	U			50.0	"	"	"	"	"
Carbon tetrachloride	U			50.0	"	"	"	"	"
1,1-Dichloropropene	U			50.0	"	"	"	"	"
Benzene	U			50.0	"	"	"	"	"
1,2-Dichloroethane	U			50.0	"	"	"	"	"
Trichloroethene	U			50.0	"	"	"	"	"
1,2-Dichloropropane	U			50.0	"	"	"	"	"
Dibromomethane	U			50.0	"	"	"	"	"
Bromodichloromethane	U			50.0	"	"	"	"	"
cis-1,3-Dichloropropene	U			50.0	"	"	"	"	"
4-Methyl-2-pentanone	U			125	"	"	"	"	"
trans-1,3-Dichloropropene	U			50.0	"	"	"	"	"
1,1,2-Trichloroethane	U			50.0	"	"	"	"	"
Tetrachloroethene	U			50.0	"	"	"	"	"
1,3-Dichloropropane	U			50.0	"	"	"	"	"
2-Hexanone	U			125	"	"	"	"	"
Dibromochloromethane	U			50.0	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			50.0	"	"	"	"	"
Chlorobenzene	U			50.0	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			50.0	"	"	"	"	"
Styrene	U			50.0	"	"	"	"	"
Bromoform	U			50.0	"	"	"	"	"
Isopropylbenzene	78.1			50.0	"	"	"	"	"
Bromobenzene	U			50.0	"	"	"	"	"
1,2,3-Trichloropropane	U			50.0	"	"	"	"	"
n-Propylbenzene	87.4			50.0	"	"	"	"	"
2-Chlorotoluene	U			50.0	"	"	"	"	"
1,3,5-Trimethylbenzene	161			50.0	"	"	"	"	"
4-Chlorotoluene	U			50.0	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			50.0	"	"	"	"	"
tert-Butylbenzene	U			50.0	"	"	"	"	"
1,2,4-Trimethylbenzene	588			50.0	"	"	"	"	"





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Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-MW002-201201 (2012005-05RE2)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
sec-Butylbenzene	U			50.0	ug/L	25	B20L008	Dec-04-20	Dec-05-20
1,3-Dichlorobenzene	U			50.0	"	"	"	"	"
p-Isopropyltoluene	U			50.0	"	"	"	"	"
1,4-Dichlorobenzene	U			50.0	"	"	"	"	"
1,2-Dichlorobenzene	U			50.0	"	"	"	"	"
n-Butylbenzene	U			50.0	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			50.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U			50.0	"	"	"	"	"
Hexachlorobutadiene	U			50.0	"	"	"	"	"
Naphthalene	58.5	B		50.0	"	"	"	"	"
1,2,3-Trichlorobenzene	U			50.0	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	10.0			99.5%		73-124	"	"	"
1,2-Dichloroethane-d4	10.6			105%		84-122	"	"	"
Toluene-d8	10.5			105%		88-108	"	"	"
4-Bromofluorobenzene	9.43			94.3%		84-108	"	"	"

A11-TB002-201201 (2012005-06RE1)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dichlorodifluoromethane	U			2.00	ug/L	1	B20L008	Dec-04-20	Dec-04-20
Chloromethane	U			2.00	"	"	"	"	"
Vinyl chloride	U			2.00	"	"	"	"	"
Bromomethane	U			2.00	"	"	"	"	"
Chloroethane	U			2.00	"	"	"	"	"
Trichlorofluoromethane	U			2.00	"	"	"	"	"
1,1-Dichloroethene	U			2.00	"	"	"	"	"
Acetone	U			12.5	"	"	"	"	"
Carbon disulfide	U			2.00	"	"	"	"	"
Methylene chloride	U			2.00	"	"	"	"	"
trans-1,2-Dichloroethene	U			2.00	"	"	"	"	"
1,1-Dichloroethane	U			2.00	"	"	"	"	"
2,2-Dichloropropane	U			2.00	"	"	"	"	"
cis-1,2-Dichloroethene	U			2.00	"	"	"	"	"
2-Butanone	U			12.5	"	"	"	"	"
Bromochloromethane	U			2.00	"	"	"	"	"



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Volatiles by GC/MS, EPA 8260C (modified)  
US EPA Region 5 LSASD Analytical Services Branch

A11-TB002-201201 (2012005-06RE1)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Chloroform	U			2.00	ug/L	1	B20L008	Dec-04-20	Dec-04-20
1,1,1-Trichloroethane	U			2.00	"	"	"	"	"
Carbon tetrachloride	U			2.00	"	"	"	"	"
1,1-Dichloropropene	U			2.00	"	"	"	"	"
Benzene	U			2.00	"	"	"	"	"
1,2-Dichloroethane	U			2.00	"	"	"	"	"
Trichloroethene	U			2.00	"	"	"	"	"
1,2-Dichloropropane	U			2.00	"	"	"	"	"
Dibromomethane	U			2.00	"	"	"	"	"
Bromodichloromethane	U			2.00	"	"	"	"	"
cis-1,3-Dichloropropene	U			2.00	"	"	"	"	"
4-Methyl-2-pentanone	U			5.00	"	"	"	"	"
Toluene	U			2.00	"	"	"	"	"
trans-1,3-Dichloropropene	U			2.00	"	"	"	"	"
1,1,2-Trichloroethane	U			2.00	"	"	"	"	"
Tetrachloroethene	U			2.00	"	"	"	"	"
1,3-Dichloropropane	U			2.00	"	"	"	"	"
2-Hexanone	U			5.00	"	"	"	"	"
Dibromochloromethane	U			2.00	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			2.00	"	"	"	"	"
Chlorobenzene	U			2.00	"	"	"	"	"
1,1,1,2-Tetrachloroethane	U			2.00	"	"	"	"	"
Ethylbenzene	U			2.00	"	"	"	"	"
m+p-Xylene	U			4.00	"	"	"	"	"
o-Xylene	U			2.00	"	"	"	"	"
Styrene	U			2.00	"	"	"	"	"
Bromoform	U			2.00	"	"	"	"	"
Isopropylbenzene	U			2.00	"	"	"	"	"
Bromobenzene	U			2.00	"	"	"	"	"
1,2,3-Trichloropropane	U			2.00	"	"	"	"	"
n-Propylbenzene	U			2.00	"	"	"	"	"
2-Chlorotoluene	U			2.00	"	"	"	"	"
1,3,5-Trimethylbenzene	U			2.00	"	"	"	"	"
4-Chlorotoluene	U			2.00	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			2.00	"	"	"	"	"
tert-Butylbenzene	U			2.00	"	"	"	"	"
1,2,4-Trimethylbenzene	U			2.00	"	"	"	"	"





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A11-TB002-201201 (2012005-06RE1)

Matrix: Water

Sampled: Dec-02-20 13:10

Received: Dec-03-20 10:57

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
sec-Butylbenzene	U			2.00	ug/L	1	B20L008	Dec-04-20	Dec-04-20
1,3-Dichlorobenzene	U			2.00	"	"	"	"	"
p-Isopropyltoluene	U			2.00	"	"	"	"	"
1,4-Dichlorobenzene	U			2.00	"	"	"	"	"
1,2-Dichlorobenzene	U			2.00	"	"	"	"	"
n-Butylbenzene	U			2.00	"	"	"	"	"
1,2-Dibromo-3-chloropropane	U			2.00	"	"	"	"	"
1,2,4-Trichlorobenzene	U			2.00	"	"	"	"	"
Hexachlorobutadiene	U			2.00	"	"	"	"	"
Naphthalene	U			2.00	"	"	"	"	"
1,2,3-Trichlorobenzene	U			2.00	"	"	"	"	"
Surrogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Dibromofluoromethane	9.82			97.7%		73-124	"	"	"
1,2-Dichloroethane-d4	10.4			103%		84-122	"	"	"
Toluene-d8	9.92			99.2%		88-108	"	"	"
4-Bromofluorobenzene	9.37			93.7%		84-108	"	"	"

**Southeast Rockford Area 11 - Groundwater Samples  
Data Validation Report**

**Sample Delivery Group (SDG) Number:** 680-192276

**Laboratory:** Eurofins Test America

**Matrix:** Groundwater

**Collection date:** 12/1/2020 & 12/2/2020

**Analysis/Methods:** Dissolved Gases - Methane - RSK-175

**Samples in SDG:**

<u>Lab ID</u>	<u>Sample Number</u>	<u>Lab ID</u>	<u>Sample Number</u>
680-192276-1	A11-MW006-201201	680-192276-7	A11-MW003-201202
680-192276-2	A11-MW130A-201201	680-192276-8	A11-MW007-201202
680-192276-3	A11-MW005-201201	680-192276-9	A11-MW007-201202-D
680-192276-4	A11-MW001-201201	680-192276-10	A11-MW002-201202
680-192276-5	A11-MW004B-201201	680-192276-11	A11-MW004A-201202
680-192276-6	A11-FB01-201201	680-192276-12	A11-TB01-201201

Data validation was performed in accordance with the specific analytical methods and the National Functional Guidelines for Organic Superfund Methods Data Review (EPA January 2017).

**Methane (RSK-175)**

<b>Precision:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Are the field duplicate relative percent differences (RPD) ≤30% (aqueous)?	Yes		
Were the Matrix Spike Duplicate RPDs ≤ 20%? (Or lab defined limits)	Yes		
Laboratory Control Spike Duplicates RPD within limits?	Yes		
Laboratory Duplicate RPDs within limits?			N/A
<u>Comments (note deviations):</u>			

<u>Field Duplicates</u>	<u>Sample</u> A11-MW007-201202	<u>Duplicate</u> A11-MW007-201202-D	<u>%RPD</u> Acceptable	<u>Qualifiers</u>	<u>Associated Samples</u>
<u>MS/MSD</u> MS/MSD 680-192276-3	<u>%RPD</u> Acceptable	<u>Limit</u>		<u>Qualifiers</u>	<u>Associated Samples</u>
<u>LCS/LCSD</u> LCS 680-647488/ 6 / 7 LCS 680-647488/ 3 / 4	<u>%RPD</u> Acceptable Acceptable	<u>Limits</u>		<u>Qualifiers</u>	<u>Associated Samples</u>
<u>Laboratory Duplicate</u> N/A	<u>%RPD</u>	<u>Limits</u>		<u>Qualifiers</u>	<u>Associated Samples</u>

<b>Accuracy:</b>	<b>Yes</b>	<b>No</b>	<b>N/A</b>
Was the Matrix Spike/Matrix Spike Duplicate criteria met? (frequency ≥ 5% and laboratory determined control limits)	Yes		
Laboratory Control Sample criteria met?	Yes		
Were the Laboratory Method Blank results all < RL?	Yes		
Were the Field Blanks results all < RL?	No		
Was the ICAL criteria met?	Yes		
Was the CCV criteria met?	Yes		
Was the Tuning criteria met?	N/A		
Were the Surrogate % recoveries within laboratory determined control limits?	N/A		
Were the Internal Standard areas within ± 50 - 150%?	N/A		
<u>Comments (note deviations):</u>			

<u>Blanks</u>	<u>Concentration</u> (mg/L)	<u>MDL /PQL</u>	<u>Qualifiers</u>	<u>Associated Samples</u>
MB 680-647488/ 8	Nondetect			



<b>Field Blank</b>		<b><u>Concentration</u></b>	<b><u>MDL /PQL</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
A11-TB01-201201	Methane	0.47 J	0.29 / 0.58	U-RL	680-192276-3	
A11-FB01-201201	Methane	0.52 J	0.29 / 0.58	U-RL	680-192276-3	
<b>Surrogates</b>		<b><u>%R</u></b>	<b><u>Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
N/A						
<b>MS/MSD</b>		<b><u>%R</u></b>	<b><u>Limits (%)</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
MS/MSD 680-192276-3		Acceptable				
<b>LCS/LCSD</b>		<b><u>%R</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
LCS 680-647488/ 6 / 7		Acceptable				
LCS 680-647488/ 3 / 4		Acceptable				
<b>ICAL</b>		<b><u>RRF</u></b>	<b><u>%RSD</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
2/17/2020 8:45		Acceptable	Acceptable			
3/04/2020 9:12		Acceptable	Acceptable			
<b>ICV / CCV</b>		<b><u>RRF</u></b>	<b><u>%D</u></b>	<b><u>Limits</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>
ICV						
3/04/2020 11:29		Acceptable	Acceptable			
CCV						
12/08/2020 17:37		Acceptable	Acceptable			
12/08/2020 17:12		Acceptable	Acceptable			
12/08/2020 20:20		Acceptable	Acceptable			
<b>Tune</b>						
N/A						
<b>Internal Standards</b>		<b><u>Area</u></b>	<b><u>Area Lower / Upper Limit</u></b>	<b><u>Qualifiers</u></b>	<b><u>Associated Samples</u></b>	
N/A						
<b>Representativeness:</b>						<b><u>Yes</u></b> <b><u>No</u></b> <b><u>N/A</u></b>
Were sampling procedures and design criteria met?						<b>Yes</b>
Were holding times met?						<b>Yes</b>
Was preservation criteria met? (0° C - 6° C)						<b>Yes</b>
Were Chain-of-Custody records complete and provided in data package?						<b>Yes</b>
<u>Comments (note deviations):</u> The cooler temperature was 4.1° C.						
<b>Preservation</b>		<b><u>Cooler Temperature (Degrees C)</u></b>	<b><u>Preservation Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>	
		Acceptable				
<b>Holding Times</b>	<b><u>Analyte</u></b>	<b><u>Days to Extraction</u></b>	<b><u>HT Criteria</u></b>	<b><u>Qualifier</u></b>	<b><u>Associated Samples</u></b>	
		Acceptable				
<b>Comparability:</b>						<b><u>Yes</u></b> <b><u>No</u></b> <b><u>N/A</u></b>
Were analytical procedures and methods followed as defined in the QAPP or field change documentation?						<b>Yes</b>
<u>Comments (note deviations):</u>						
<b>Completeness (90%):</b>						<b><u>Yes</u></b> <b><u>No</u></b> <b><u>N/A</u></b>
Are all data in this SDG usable?						<b>Yes</b>
<u>Comments (note deviations):</u>						

**Sensitivity:**

Are MDLs present and reported?  
Do the reporting limits meet project requirements?

Comments (note deviations):

Yes	No	N/A
Yes		
Yes		

**Comment:**

As stated in the case narrative, the MS/MSD was spiked at the concentration range meant for the TCD detector. The methane results for the FID detector were over the calibration range as a result of the error. The recovery areas are within limits for both FID and TCD detectors. Both the FID and TCD detectors are being reported for the MS/MSD.

Data is usable with appropriate qualifiers applied.

Data Validator:	<u>Kristine Molloy</u>
Data Reviewer:	<u>Cherie Zakowski</u>

Date:	<u>5/2/2021</u>
Date:	<u>5/4/2021</u>



# Detection Summary

Client: CDM Smith, Inc.  
Project/Site: Methane Analysis - SE Rockford Area 11

Job ID: 680-192276-1

## Client Sample ID: A11-MW006-201201

## Lab Sample ID: 680-192276-1

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	8100		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW130A-201201

## Lab Sample ID: 680-192276-2

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	1.4		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW005-201201

## Lab Sample ID: 680-192276-3

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.48	J	0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW001-201201

## Lab Sample ID: 680-192276-4

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	11		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW004B-201201

## Lab Sample ID: 680-192276-5

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	20		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-FB01-201201

## Lab Sample ID: 680-192276-6

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.52	J	0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW003-201202

## Lab Sample ID: 680-192276-7

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	6600		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW007-201202

## Lab Sample ID: 680-192276-8

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	31000		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW007-201202-D

## Lab Sample ID: 680-192276-9

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	29000		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW002-201202

## Lab Sample ID: 680-192276-10

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane (TCD)	31000		390	39	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-MW004A-201202

## Lab Sample ID: 680-192276-11

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	250		0.58	0.29	ug/L	1			RSK-175	Total/NA

## Client Sample ID: A11-TB01-201201

## Lab Sample ID: 680-192276-12

Analyte	Result	Qualifier	RL	MDL	Unit	Dil	Fac	D	Method	Prep Type
Methane	0.47	J	0.58	0.29	ug/L	1			RSK-175	Total/NA

This Detection Summary does not include radiochemical test results.

Eurofins TestAmerica, Savannah

# Client Sample Results

Client: CDM Smith, Inc.  
Project/Site: Methane Analysis - SE Rockford Area 11

Job ID: 680-192276-1

**Client Sample ID: A11-MW006-201201**

**Lab Sample ID: 680-192276-1**

Date Collected: 12/01/20 12:11

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	8100		390	39	ug/L			12/08/20 16:46	1

**Client Sample ID: A11-MW130A-201201**

**Lab Sample ID: 680-192276-2**

Date Collected: 12/01/20 09:35

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	1.4		0.58	0.29	ug/L			12/08/20 16:59	1

**Client Sample ID: A11-MW005-201201**

**Lab Sample ID: 680-192276-3**

Date Collected: 12/01/20 13:50

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.48	J	0.58	0.29	ug/L			12/08/20 17:45	1

**Client Sample ID: A11-MW001-201201**

**Lab Sample ID: 680-192276-4**

Date Collected: 12/01/20 15:31

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	11		0.58	0.29	ug/L			12/08/20 18:24	1

**Client Sample ID: A11-MW004B-201201**

**Lab Sample ID: 680-192276-5**

Date Collected: 12/01/20 16:45

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	20		0.58	0.29	ug/L			12/08/20 18:36	1

**Client Sample ID: A11-FB01-201201**

**Lab Sample ID: 680-192276-6**

Date Collected: 12/01/20 17:05

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.52	J	0.58	0.29	ug/L			12/08/20 18:49	1

**Client Sample ID: A11-MW003-201202**

**Lab Sample ID: 680-192276-7**

Date Collected: 12/02/20 08:45

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	6600		390	39	ug/L			12/08/20 19:02	1



# Client Sample Results

Client: CDM Smith, Inc.

Job ID: 680-192276-1

Project/Site: Methane Analysis - SE Rockford Area 11

**Client Sample ID: A11-MW007-201202**

**Lab Sample ID: 680-192276-8**

Date Collected: 12/02/20 10:50

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	31000		390	39	ug/L			12/08/20 19:15	1

**Client Sample ID: A11-MW007-201202-D**

**Lab Sample ID: 680-192276-9**

Date Collected: 12/02/20 10:50

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	29000		390	39	ug/L			12/08/20 19:28	1

**Client Sample ID: A11-MW002-201202**

**Lab Sample ID: 680-192276-10**

Date Collected: 12/02/20 13:10

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane (TCD)	31000		390	39	ug/L			12/08/20 19:41	1

**Client Sample ID: A11-MW004A-201202**

**Lab Sample ID: 680-192276-11**

Date Collected: 12/02/20 15:00

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	250		0.58	0.29	ug/L			12/08/20 19:54	1

**Client Sample ID: A11-TB01-201201**

**Lab Sample ID: 680-192276-12**

Date Collected: 12/01/20 08:00

Matrix: Water

Date Received: 12/04/20 11:00

**Method: RSK-175 - Dissolved Gases (GC)**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methane	0.47	J	0.58	0.29	ug/L			12/08/20 20:07	1

## Default Detection Limits

Client: CDM Smith, Inc.

Job ID: 680-192276-1

Project/Site: Methane Analysis - SE Rockford Area 11

### Method: RSK-175 - Dissolved Gases (GC)

Analyte	RL	MDL	Units
Methane	0.58	0.29	ug/L
Methane (TCD)	390	39	ug/L